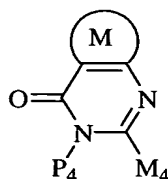


WHAT IS CLAIMED IS:

1. A compound of formula I:



I

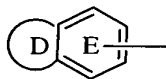
or a stereoisomer or pharmaceutically acceptable salt thereof, wherein;

ring M is a 5 or 6 membered aromatic or dihydro-aromatic ring consisting of: carbon atoms and 0-3 heteroatoms selected from O, S(O)_p, and N;

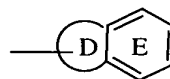
ring M is substituted with 0-3 R^{1a} and 0-1 carbonyl groups;

one of P₄ and M₄ is -Z-A-B and the other -G₁-G;

G is a group of formula IIa or IIb:



IIa



IIb

ring D, including the two atoms of Ring E to which it is attached, is a 5-6 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and S(O)_p;

ring D is substituted with 0-2 R and has 0-3 ring double bonds;

E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, and pyridazinyl, and is substituted with 1-2 R;

alternatively, ring D is absent and ring E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, pyridazinyl, pyrrolyl, pyrazolyl, imidazolyl, isoxazolyl, oxazolyl, triazolyl, thienyl, and thiazolyl, and ring E is substituted with 1-2 R;

alternatively, ring D is absent and ring E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, pyridazinyl, pyrrolyl, pyrazolyl, imidazolyl, isoxazolyl, oxazolyl, triazolyl, thienyl, and thiazolyl, and ring E is substituted with 1 R and with a 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, wherein the heterocycle is substituted with 0-1 carbonyls, 1-2 R, and 0-3 ring double bonds;

R is selected from H, C₁₋₄ alkyl, F, Cl, Br, I, OH, OCH₃, -OCH₂CH₃,

-OCH(CH₃)₂, -OCH₂CH₂CH₃, CN, -C(=NR⁸)NR⁷R⁹, -NHC(=NR⁸)NR⁷R⁹,
 -NR⁸CH(=NR⁷), NH₂, -NH(C₁₋₃ alkyl), -N(C₁₋₃ alkyl)₂, -C(=NH)NH₂, -CH₂NH₂,
 -CH₂NH(C₁₋₃ alkyl), -CH₂N(C₁₋₃ alkyl)₂, -CH₂CH₂NH₂, -CH₂CH₂NH(C₁₋₃ alkyl),
 -CH₂CH₂N(C₁₋₃ alkyl)₂, -(CR⁸R⁹)_tC(O)H, -(CR⁸R⁹)_tC(O)R^{2c}, -(CR⁸R⁹)_tNR⁷R⁸,
 5 -(CR⁸R⁹)_tC(O)NR⁷R⁸, -(CR⁸R⁹)_tNR⁷C(O)R⁷, -(CR⁸R⁹)_tOR³,
 -(CR⁸R⁹)_tS(O)_pNR⁷R⁸, -(CR⁸R⁹)_tNR⁷S(O)_pR⁷, -(CR⁸R⁹)_tSR³, -(CR⁸R⁹)_tS(O)R³,
 -(CR⁸R⁹)_tS(O)₂R³, and -OCF₃, provided that S(O)_pR⁷ forms other than S(O)₂H or
 S(O)H;

alternatively, when 2 R groups are attached to adjacent atoms, they combine to
 10 form methylenedioxy or ethylenedioxy;

A is selected from: C₃₋₁₀ carbocycle substituted with 0-2 R⁴, and 5-12
 membered heterocycle substituted with 0-2 R⁴ and consisting of: carbon atoms and
 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

B is selected from: Y, X-Y, -(CH₂)₀₋₂C(O)NR²R^{2a}, -(CH₂)₀₋₂NR²R^{2a},
 15 -C(=NR²)NR²R^{2a}, and -NR²C(=NR²)NR²R^{2a}, provided that Z and B are attached to
 different atoms on A;

X is selected from -(CR²R^{2a})₁₋₄-, -CR²(CR²R^{2b})(CH₂)_t-, -C(O)-, -C(=NR^{1b})-,
 -CR²(NR^{1b}R²)-, -CR²(OR²)-, -CR²(SR²)-, -C(O)CR²R^{2a}-, -CR²R^{2a}C(O)-, -S-, -S(O)-,
 -S(O)₂-, -SCR²R^{2a}-, -S(O)CR²R^{2a}-, -S(O)₂CR²R^{2a}-, -CR²R^{2a}S-, -CR²R^{2a}S(O)-,
 20 -CR²R^{2a}S(O)₂-, -S(O)₂NR²-, -NR²S(O)₂-, -NR²S(O)₂CR²R^{2a}-, -CR²R^{2a}S(O)₂NR²-,
 -NR²S(O)₂NR²-, -C(O)NR²-, -NR²C(O)-, -C(O)NR²CR²R^{2a}-, -NR²C(O)CR²R^{2a}-,
 -CR²R^{2a}C(O)NR²-, -CR²R^{2a}NR²C(O)-, -NR²C(O)O-, -OC(O)NR²-, -NR²C(O)NR²-,
 -NR²-, -NR²CR²R^{2a}-, -CR²R^{2a}NR²-, O, -CR²R^{2a}O-, and -OCR²R^{2a}-;

Y is selected from: C₃₋₁₀ carbocycle substituted with 0-2 R^{4a}, and 5-10
 25 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from
 the group consisting of N, O, and S(O)_p substituted with 0-2 R^{4a};

G₁ is absent or is selected from -(CR³R^{3a})₁₋₅-,
 -(CR³R^{3a})₀₋₂CR³=CR³(CR³R^{3a})₀₋₂-, -(CR³R^{3a})₀₋₂C≡C(CR³R^{3a})₀₋₂-,
 -(CR³R^{3a})_uC(O)(CR³R^{3a})_w-, -(CR³R^{3a})_uC(O)O(CR³R^{3a})_w-,
 30 -(CR³R^{3a})_uOC(O)(CR³R^{3a})_w-, -(CR³R^{3a})_uO(CR³R^{3a})_w-;

- CR³R^{3a})_uNR^{3b}(CR³R^{3a})_{w-}, -(CR³R^{3a})_uC(O)NR^{3b}(CR³R^{3a})_{w-},
-(CR³R^{3a})_uNR^{3b}C(O)(CR³R^{3a})_{w-}, -(CR³R^{3a})_uOC(O)NR^{3b}(CR³R^{3a})_{w-},
-(CR³R^{3a})_uNR^{3b}C(O)O(CR³R^{3a})_{w-}, -(CR³R^{3a})_uNR^{3b}C(O)NR^{3b}(CR³R^{3a})_{w-},
-(CR³R^{3a})_uNR^{3b}C(S)NR^{3b}(CR³R^{3a})_{w-}, -(CR³R^{3a})_uS(CR³R^{3a})_{w-},
5 -(CR³R^{3a})_uS(O)(CR³R^{3a})_{w-}, -(CR³R^{3a})_uS(O)₂(CR³R^{3a})_{w-},
-(CR³R^{3a})_uS(O)NR^{3b}(CR³R^{3a})_{w-}, -(CR³R^{3a})_uNR^{3b}S(O)₂(CR³R^{3a})_{w-},
-(CR³R^{3a})_uS(O)₂NR^{3b}(CR³R^{3a})_{w-}, -(CR³R^{3a})_uNR^{3b}S(O)₂NR^{3b}(CR³R^{3a})_{w-},
-(CR³R^{3a})_uNR^{3e}(CR³R^{3a})_{w-}, -(CR³R^{3a})_uC(O)(CR³R^{3a})_uC(O)(CR³R^{3a})_{w-},
-(CR³R^{3a})_uNR^{3b}(CR³R^{3a})_uC(O)NR^{3b}(CR³R^{3a})_{w-},
10 -(CR³R^{3a})_uNR^{3b}C(O)(CR³R^{3a})_uC(O)(CR³R^{3a})_{w-},
-(CR³R^{3a})_uC(O)(CR³R^{3a})_uC(O)NR^{3b}(CR³R^{3a})_{w-},
-(CR³R^{3a})_uNR^{3b}C(O)(CR³R^{3a})_uC(O)NR^{3b}(CR³R^{3a})_{w-},
-(CR³R^{3a})_uS(O)NR^{3b}C(O)(CR³R^{3a})_{w-}, -(CR³R^{3a})_uC(O)NR^{3b}S(O)₂(CR³R^{3a})_{w-}, and
-(CR³R^{3a})_uS(O)₂NR^{3b}C(O)NR^{3b}CR³R^{3a})_{w-}, wherein u + w total 0, 1, 2, 3, or 4,
15 provided that G₁ does not form a N-S, NCH₂N, NCH₂O, or NCH₂S bond with either
group to which it is attached;
Z is selected from a bond, -(CR³R^{3a})_{l-4-}, -(CR³R^{3a})_qO(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qNR^{3b}(CR³R^{3a})_{q1-}, -(CR³R^{3a})_qC(O)(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qC(O)O(CR³R^{3a})_{q1-}, -(CR³R^{3a})_qOC(O)(CR³R^{3a})_{q1-},
20 -(CR³R^{3a})_qC(O)NR^{3b}(CR³R^{3a})_{q1-}, -(CR³R^{3a})_qNR^{3b}C(O)(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qOC(O)O(CR³R^{3a})_{q1-}, -(CR³R^{3a})_qOC(O)NR^{3b}(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qNR^{3b}C(O)O(CR³R^{3a})_{q1-}, -(CR³R^{3a})_qNR^{3b}C(O)NR^{3b}(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qC(O)(CR³R^{3a})_qC(O)(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qNR^{3b}(CR³R^{3a})_qC(O)NR^{3b}(CR³R^{3a})_{q1-},
25 -(CR³R^{3a})_qNR^{3b}C(O)(CR³R^{3a})_qC(O)(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qC(O)(CR³R^{3a})_qC(O)NR^{3b}(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qNR^{3b}C(O)(CR³R^{3a})_qC(O)NR^{3b}(CR³R^{3a})_{q1-}, -(CR³R^{3a})_qS(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qS(O)(CR³R^{3a})_{q1-}, -(CR³R^{3a})_qS(O)₂(CR³R^{3a})_{q1-},
-(CR³R^{3a})_qSO₂NR^{3b}(CR³R^{3a})_{q1-}, -(CR³R^{3a})_qNR^{3b}SO₂(CR³R^{3a})_{q1-},

$-(\text{CR}^3\text{R}^{3a})_q\text{S}(\text{O})\text{NR}^{3b}\text{C}(\text{O})(\text{CR}^3\text{R}^{3a})_{q1}-$, $-(\text{CR}^3\text{R}^{3a})_q\text{C}(\text{O})\text{NR}^{3b}\text{S}(\text{O})_2(\text{CR}^3\text{R}^{3a})_{q1}-$,
and $-(\text{CR}^3\text{R}^{3a})_q\text{NR}^{3b}\text{SO}_2\text{NR}^{3b}(\text{CR}^3\text{R}^{3a})_{q1}-$, wherein $q + q1$ total 0, 1, 2, 3, or 4,
provided that Z does not form a N-S, NCH_2N , NCH_2O , or NCH_2S bond with either
group to which it is attached;

- 5 R^{1a} , at each occurrence, is selected from H, $-(\text{CR}^3\text{R}^{3a})_r\text{R}^{1b}$,
 $-(\text{CR}^3\text{R}^{3a})_r\text{CR}^3\text{R}^{1b}\text{R}^{1b}$, $-(\text{CR}^3\text{R}^{3a})_r\text{O}-(\text{CR}^3\text{R}^{3a})_r\text{R}^{1b}$,
 $-(\text{CR}^3\text{R}^{3a})_r\text{NR}^2-(\text{CR}^3\text{R}^{3a})_r\text{R}^{1b}$, $-(\text{CR}^3\text{R}^{3a})_r\text{S}(\text{O})_p-(\text{CR}^3\text{R}^{3a})_r\text{R}^{1b}$,
 $-(\text{CR}^3\text{R}^{3a})_r\text{CO}_2-(\text{CR}^3\text{R}^{3a})_r\text{R}^{1b}$, $-(\text{CR}^3\text{R}^{3a})_r\text{C}(\text{O})\text{NR}^2-(\text{CR}^3\text{R}^{3a})_r\text{R}^{1b}$,
 $-(\text{CR}^3\text{R}^{3a})_r\text{C}(\text{O})-(\text{CR}^3\text{R}^{3a})_r\text{R}^{1b}$, C_{2-6} alkenylene- R^{1b} , C_{2-6} alkynylene- R^{1b} , and
10 $-(\text{CR}^3\text{R}^{3a})_r\text{C}(=\text{NR}^{1b})\text{NR}^{3b}\text{R}^{1b}$, provided that R^{1a} forms other than an N-halo, N-S,
O-O, or N-CN bond;

- alternatively, when two R^{1a} groups are attached to adjacent atoms, together
with the atoms to which they are attached, they form a 5-7 membered ring consisting
of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and
15 $\text{S}(\text{O})_p$, this ring being substituted with 0-2 R^{4b} and 0-3 ring double bonds;

- R^{1b} is selected from H, C_{1-3} alkyl, F, Cl, Br, I, CN, NO_2 , CHO, $-(\text{CF}_2)_r\text{CF}_3$,
 $-(\text{CR}^3\text{R}^{3a})_r\text{OR}^2$, $-\text{NR}^2\text{R}^{2a}$, $-\text{C}(\text{O})\text{R}^{2b}$, $-\text{CO}_2\text{R}^{2b}$, $-\text{OC}(\text{O})\text{R}^2$, $-\text{CH}(\text{CH}_2\text{OR}^2)_2$,
 $-(\text{CF}_2)_r\text{CO}_2\text{R}^{2a}$, $-\text{S}(\text{O})_p\text{R}^{2b}$, $-\text{NR}^2(\text{CH}_2)_r\text{OR}^2$, $-\text{C}(=\text{NR}^{2c})\text{NR}^2\text{R}^{2a}$, $-\text{NR}^2\text{C}(\text{O})\text{R}^{2b}$,
 $-\text{NR}^2\text{C}(\text{O})\text{NR}^2\text{R}^{2a}$, $-\text{NR}^2\text{C}(\text{O})_2\text{R}^{2a}$, $-\text{OC}(\text{O})\text{NR}^2\text{R}^{2a}$, $-\text{C}(\text{O})\text{NR}^2\text{R}^{2a}$,
20 $-\text{C}(\text{O})\text{NR}^2(\text{CH}_2)_r\text{OR}^2$, $-\text{SO}_2\text{NR}^2\text{R}^{2a}$, $-\text{NR}^2\text{SO}_2\text{R}^2$, $-\text{C}(\text{O})\text{NR}^2\text{SO}_2\text{R}^2$, C_{3-6} carbocycle
substituted with 0-2 R^{4b} , and 5-10 membered heterocycle substituted with 0-2 R^{4b}
and consisting of carbon atoms and from 1-4 heteroatoms selected from the group
consisting of N, O, and $\text{S}(\text{O})_p$, provided that R^{1b} forms other than an O-O, N-halo,
N-S, or N-CN bond and provided that $\text{S}(\text{O})_p\text{R}^2$ forms other than $\text{S}(\text{O})_2\text{H}$ or $\text{S}(\text{O})\text{H}$;

- 25 R^2 , at each occurrence, is selected from H, CF_3 , C_{1-6} alkyl, benzyl,
 $-(\text{CH}_2)_r\text{C}_{3-10}$ carbocycle substituted with 0-2 R^{4b} , and $-(\text{CH}_2)_r\text{5-10}$ membered
heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4
heteroatoms selected from the group consisting of N, O, and $\text{S}(\text{O})_p$;

R^{2a} , at each occurrence, is selected from H, CF_3 , C_{1-6} alkyl, benzyl,

-(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-2 R^{4b}, and -(CH₂)_r-5-10 membered heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

alternatively, R² and R^{2a}, together with the nitrogen atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-2 R^{4b} and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{2b}, at each occurrence, is selected from CF₃, C₁₋₄ alkoxy, C₁₋₆ alkyl substituted with 0-2 R^{4b}, -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-2 R^{4b}, and -(CH₂)_r-5-10 membered heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{2c}, at each occurrence, is selected from CF₃, OH, C₁₋₄ alkoxy, C₁₋₆ alkyl, -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-2 R^{4b}, and -(CH₂)_r-5-10 membered heterocycle substituted with 0-2 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R³, at each occurrence, is selected from H, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂, CH(CH₃)CH₂CH₃, C(CH₃)₃, benzyl, and phenyl;

R^{3a}, at each occurrence, is selected from H, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂, CH(CH₃)CH₂CH₃, C(CH₃)₃, benzyl, and phenyl;

R^{3b}, at each occurrence, is selected from H, C₁₋₆ alkyl substituted with 0-2 R^{1a}, C₂₋₆ alkenyl substituted with 0-2 R^{1a}, C₂₋₆ alkynyl substituted with 0-2 R^{1a}, -(C₀₋₄ alkyl)-5-10 membered carbocycle substituted with 0-3 R^{1a}, and -(C₀₋₄ alkyl)-5-10 membered heterocycle substituted with 0-3 R^{1a} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{3c} , at each occurrence, is selected from CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, and phenyl;

R^{3d} , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$,
5 $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, C_{1-4} alkyl-phenyl, and $C(=O)R^{3c}$;

R^{3e} , is selected from H, $-S(O)_2NHR^3$, $-C(O)R^3$, $-C(O)NHR^3$, $-C(O)OR^{3f}$, $-S(O)R^{3f}$, $-S(O)_2R^{3f}$, C_{1-6} alkyl substituted with 0-2 R^{1a} , C_{2-6} alkenyl substituted with 0-2 R^{1a} , C_{2-6} alkynyl substituted with 0-2 R^{1a} , $-(C_{0-4}$ alkyl)-5-10 membered carbocycle substituted with 0-3 R^{1a} , and $-(C_{0-4}$ alkyl)-5-10 membered heterocycle substituted with 0-3 R^{1a} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{3f} , at each occurrence, is selected from: C_{1-6} alkyl substituted with 0-2 R^{1a} , C_{2-6} alkenyl substituted with 0-2 R^{1a} , C_{2-6} alkynyl substituted with 0-2 R^{1a} ,
15 $-(C_{0-4}$ alkyl)-5-10 membered carbocycle substituted with 0-3 R^{1a} , and $-(C_{0-4}$ alkyl)-5-10 membered heterocycle substituted with 0-3 R^{1a} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^4 , at each occurrence, is selected from H, =O, $(CR^3R^{3a})_rOR^2$, F, Cl, Br, I, C_{1-4} alkyl, $-(CR^3R^{3a})_rCN$, $-(CR^3R^{3a})_rNO_2$, $-(CR^3R^{3a})_rNR^2R^{2a}$, $-(CR^3R^{3a})_rC(O)R^{2c}$,
20 $-(CR^3R^{3a})_rNR^2C(O)R^{2b}$, $-(CR^3R^{3a})_rC(O)NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2C(O)NR^2R^{2a}$, $-(CR^3R^{3a})_rC(=NR^2)NR^2R^{2a}$, $-(CR^3R^{3a})_rC(=NS(O)_2R^{5a})NR^2R^{2a}$, $-(CR^3R^{3a})_rNHC(=NR^2)NR^2R^{2a}$, $-(CR^3R^{3a})_rC(O)NHC(=NR^2)NR^2R^{2a}$, $-(CR^3R^{3a})_rSO_2NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2SO_2NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2SO_2-C_{1-4}$ alkyl, $-(CR^3R^{3a})_rNR^2SO_2R^{5a}$, $-(CR^3R^{3a})_rS(O)_pR^{5a}$, $-(CR^3R^{3a})_r(CF_2)_rCF_3$, $-NHCH_2R^{1b}$,
25 $-OCH_2R^{1b}$, $-SCH_2R^{1b}$, $-N(CH_2)_2(CH_2)_tR^{1b}$, $-O(CH_2)_2(CH_2)_tR^{1b}$, $-S(CH_2)_2(CH_2)_tR^{1b}$, $-(CR^3R^{3a})_r$ -5-6 membered carbocycle substituted with 0-1 R^5 , and a $-(CR^3R^{3a})_r$ -5-6 membered heterocycle substituted with 0-1 R^5 and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

- R^{4a} , at each occurrence, is selected from H, =O, $-(CR^3R^{3a})_rOR^2$, $-(CR^3R^{3a})_rF$, $-(CR^3R^{3a})_rBr$, $-(CR^3R^{3a})_rCl$, $-(CR^3R^{3a})_rI$, C_{1-4} alkyl, $-(CR^3R^{3a})_rCN$, $-(CR^3R^{3a})_rNO_2$, $-(CR^3R^{3a})_rNR^2R^{2a}$, $-(CR^3R^{3a})_rC(O)R^{2c}$, $-(CR^3R^{3a})_rNR^2C(O)R^{2b}$, $-(CR^3R^{3a})_rC(O)NR^2R^{2a}$, $-(CR^3R^{3a})_rN=CHOR^3$,
 5 $-(CR^3R^{3a})_rC(O)NH(CH_2)_2NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2C(O)NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2C(O)OR^2$, $-(CR^3R^{3a})_rC(=NR^2)NR^2R^{2a}$, $-(CR^3R^{3a})_rNHC(=NR^2)NR^2R^{2a}$, $-(CR^3R^{3a})_rSO_2NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2SO_2NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2SO_2-C_{1-4}$ alkyl, $-(CR^3R^{3a})_rC(O)NHSO_2-C_{1-4}$ alkyl, $-(CR^3R^{3a})_rNR^2SO_2R^5$, $-(CR^3R^{3a})_rS(O)_pR^5$,
 10 $-(CR^3R^{3a})_r(CF_2)_rCF_3$, $-(CR^3R^{3a})_{r-3-10}$ membered carbocycle substituted with 0-1 R^5 , and a $-(CR^3R^{3a})_{r-3-10}$ membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-1 R^5 ;

- R^{4b} , at each occurrence, is selected from H, =O, $-(CR^3R^{3a})_rOR^3$,
 15 $-(CR^3R^{3a})_rF$, $-(CR^3R^{3a})_rCl$, $-(CR^3R^{3a})_rBr$, $-(CR^3R^{3a})_rI$, C_{1-4} alkyl, $-(CR^3R^{3a})_rCN$, $-(CR^3R^{3a})_rNO_2$, $-(CR^3R^{3a})_rNR^3R^{3a}$, $-(CR^3R^{3a})_rC(O)R^3$, $-(CR^3R^{3a})_rC(O)OR^{3c}$, $-(CR^3R^{3a})_rNR^3C(O)R^{3a}$, $-(CR^3R^{3a})_rC(O)NR^3R^{3a}$, $-(CR^3R^{3a})_rNR^3C(O)NR^3R^{3a}$, $-(CR^3R^{3a})_rC(=NR^3)NR^3R^{3a}$, $-(CR^3R^{3a})_rNR^3C(=NR^3)NR^3R^{3a}$, $-(CR^3R^{3a})_rSO_2NR^3R^{3a}$, $-(CR^3R^{3a})_rNR^3SO_2NR^3R^{3a}$, $-(CR^3R^{3a})_rNR^3SO_2-C_{1-4}$ alkyl,
 20 $-(CR^3R^{3a})_rNR^3SO_2CF_3$, $-(CR^3R^{3a})_rNR^3SO_2$ -phenyl, $-(CR^3R^{3a})_rS(O)_pCF_3$, $-(CR^3R^{3a})_rS(O)_p-C_{1-4}$ alkyl, $-(CR^3R^{3a})_rS(O)_p$ -phenyl, and $-(CR^3R^{3a})_r(CF_2)_rCF_3$;

- R^5 , at each occurrence, is selected from H, C_{1-6} alkyl, =O, $-(CH_2)_rOR^3$, F, Cl, Br, I, -CN, NO_2 , $-(CH_2)_rNR^3R^{3a}$, $-(CH_2)_rC(O)R^3$, $-(CH_2)_rC(O)OR^{3c}$, $NR^3C(O)R^{3a}$, $C(O)NR^3R^{3a}$, $NR^3C(O)NR^3R^{3a}$, $CH(=NOR^{3d})$, $C(=NR^3)NR^3R^{3a}$,
 25 $NR^3C(=NR^3)NR^3R^{3a}$, $SO_2NR^3R^{3a}$, $NR^3SO_2NR^3R^{3a}$, $NR^3SO_2-C_{1-4}$ alkyl, $NR^3SO_2CF_3$, NR^3SO_2 -phenyl, $S(O)_pCF_3$, $S(O)_p-C_{1-4}$ alkyl, $S(O)_p$ -phenyl, $(CF_2)_rCF_3$, phenyl substituted with 0-2 R^6 , naphthyl substituted with 0-2 R^6 , and benzyl substituted with 0-2 R^6 ;

R^{5a} , at each occurrence, is selected from C_{1-6} alkyl, $-(CH_2)_rOR^3$,

$-(CH_2)_rNR^3R^{3a}$, $-(CH_2)_rC(O)R^3$, $-(CH_2)_rC(O)OR^{3c}$, $-(CH_2)_rNR^3C(O)R^{3a}$,
 $-(CH_2)_rC(O)NR^3R^{3a}$, $-(CF_2)_rCF_3$, phenyl substituted with 0-2 R^6 , naphthyl
substituted with 0-2 R^6 , and benzyl substituted with 0-2 R^6 , provided that R^{5a} does
not form a S-N or $S(O)_p-C(O)$ bond;

5 R^6 , at each occurrence, is selected from H, OH, $-(CH_2)_rOR^2$, F, Cl, Br, I,
 C_{1-4} alkyl, -CN, NO_2 , $-(CH_2)_rNR^2R^{2a}$, $-(CH_2)_rC(O)R^{2b}$, $-NR^2C(O)R^{2b}$,
 $-NR^2C(O)NR^2R^{2a}$, $-C(=NH)NH_2$, $-NHC(=NH)NH_2$, $-SO_2NR^2R^{2a}$, $-NR^2SO_2NR^2R^{2a}$,
and $-NR^2SO_2C_{1-4}$ alkyl;

R^7 , at each occurrence, is selected from H, OH, C_{1-6} alkyl, C_{1-6} alkyl-C(O)-,
10 C_{1-6} alkyl-O-, $(CH_2)_n$ -phenyl, C_{1-4} alkyl-OC(O)-, C_{6-10} aryl-O-, C_{6-10} aryl-OC(O)-,
 C_{6-10} aryl- CH_2 -C(O)-, C_{1-4} alkyl-C(O)O- C_{1-4} alkyl-OC(O)-, C_{6-10} aryl-C(O)O- C_{1-4}
alkyl-OC(O)-, C_{1-6} alkyl- NH_2 -C(O)-, phenyl- NH_2 -C(O)-, and phenyl C_{1-4} alkyl-
C(O)-;

R^8 , at each occurrence, is selected from H, C_{1-6} alkyl, and $-(CH_2)_n$ -phenyl;
15 alternatively, R^7 and R^8 , when attached to the same nitrogen, combine to form
a 5-10 membered heterocyclic ring consisting of carbon atoms and 0-2 additional
heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^9 , at each occurrence, is selected from H, C_{1-6} alkyl, and $-(CH_2)_n$ -phenyl;
n, at each occurrence, is selected from 0, 1, 2, and 3;
20 p, at each occurrence, is selected from 0, 1, and 2;
r, at each occurrence, is selected from 0, 1, 2, 3, 4, 5, and 6; and
t, at each occurrence, is selected from 0, 1, 2, and 3;
provided that:

(a) when Z and G_1 are absent, A is phenyl or pyridyl, G is phenyl, pyridyl, or
25 thienyl, at least one R is other than a substituted or unsubstituted group
selected from amidino, guanidino, guanidine-methyl, iminoamino,
iminoamino-methyl, amino, amino-methyl, and pyridyl, then B is other than
cycloalkyl, $(CH_2)_{0-2}C(O)NR^2R^{2a}$, or $(CH_2)_{0-2}NR^2R^{2a}$, wherein substituted
includes being cyclized with an additional heteroatom being optionally
30 present;

(b) when Z and G₁ are absent, G is phenyl or pyridyl, and A is phenyl, pyridyl, furanyl, or thienyl, then B is other than a substituted or unsubstituted group selected from amidino, guanidino, guanidine-methyl, iminoamino, iminoamino-methyl, amino, amino-methyl, aminosulfonyl-phenyl, and pyridyl, wherein substituted includes being cyclized with an additional heteroatom being optionally present; and

(c) when G-G₁ is hydroxy-phenyl or alkoxy-phenyl, then B is other than acyclic or cyclic-amino-alkoxy.

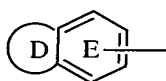
2. A compound according to Claim 1, wherein:

ring M is selected from phenyl, pyrrole, furan, thiophene, pyrazole, imidazole, isoxazole, oxazole, isothiazole, thiazole, 1,2,3-triazole, 1,2,3-oxadiazole, 1,2,3-thiadiazole, pyridine, pyrimidine, pyridazine, pyrazine, 1,2,3-triazine, 1,2,4-triazine, and 1,2,3,4-tetrazine;

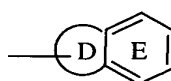
ring M is substituted with 0-3 R^{1a};

one of P₄ and M₄ is -Z-A-B and the other -G₁-G;

G is a group of formula IIa or IIb:



IIa



IIb

ring D, including the two atoms of Ring E to which it is attached, is a 5-6 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and S(O)_p;

ring D is substituted with 0-2 R and has 0-3 ring double bonds;

E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, and pyridazinyl, and is substituted with 1-2 R;

alternatively, ring D is absent, and ring E is selected from phenyl, pyridyl, pyrimidyl, and thienyl, and ring E is substituted with 1-2 R;

alternatively, ring D is absent, ring E is selected from phenyl, pyridyl, and thienyl, and ring E is substituted with a 5 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p,

wherein the heterocycle is substituted with 0-1 carbonyls, 1-2 R, and 0-3 ring double bonds;

R is selected from H, C₁₋₄ alkyl, F, Cl, OH, OCH₃, -OCH₂CH₃, -OCH(CH₃)₂, CN, -C(=NH)NH₂, NH₂, -NH(C₁₋₃ alkyl), -N(C₁₋₃ alkyl)₂, -C(=NH)NH₂, -CH₂NH₂,
 5 -CH₂NH(C₁₋₃ alkyl), -CH₂N(C₁₋₃ alkyl)₂, -(CR⁸R⁹)_tNR⁷R⁸, -C(O)NR⁷R⁸,
 -CH₂C(O)NR⁷R⁸, -S(O)_pNR⁷R⁸, -CH₂S(O)_pNR⁷R⁸, and -OCF₃;

alternatively, when 2 R groups are attached to adjacent atoms, they combine to form methylenedioxy or ethylenedioxy;

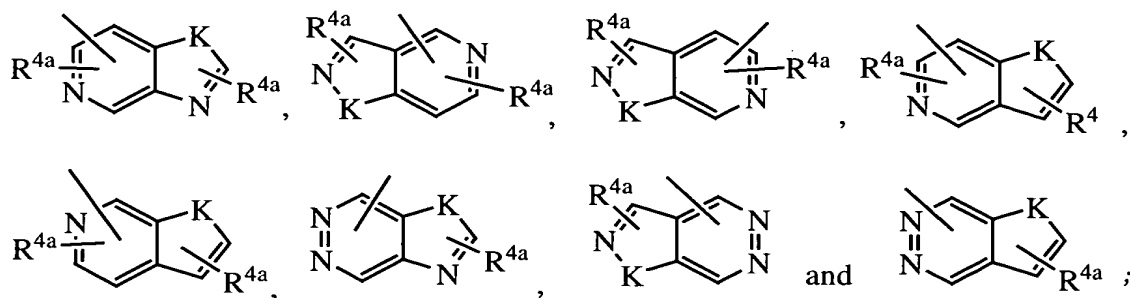
A is selected from one of the following rings and is substituted with 0-2 R⁴;
 10 phenyl, piperidinyl, piperazinyl, pyridyl, pyrimidyl, furanyl, morpholinyl, thiophenyl, pyrrolyl, pyrrolidinyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, pyrazolyl, imidazolyl, oxadiazolyl, thiadiazolyl, triazolyl, 1,2,3-oxadiazolyl, 1,2,4-oxadiazolyl, 1,2,5-oxadiazolyl, 1,3,4-oxadiazolyl, 1,2,3-thiadiazolyl, 1,2,4-thiadiazolyl, 1,2,5-thiadiazolyl, 1,3,4-thiadiazolyl, 1,2,3-triazolyl, 1,2,4-triazolyl, 1,2,5-triazolyl,
 15 1,3,4-triazolyl, benzofuranyl, benzothiofuranyl, indolyl, benzimidazolyl, benzoxazolyl, benzthiazolyl, indazolyl, benzisoxazolyl, benzisothiazolyl, and isoindazolyl;

B is selected from Y, X-Y, -CH₂NR²R^{2a}, and -CH₂CH₂NR²R^{2a};

X is selected from -(CR²R^{2a})₁₋₄-, -C(O)-, -C(=NR^{1b})-, -CR²(NR^{1b}R²)-,
 20 -C(O)CR²R^{2a}-, -CR²R^{2a}C(O)-, -C(O)NR²-, -NR²C(O)-, -C(O)NR²CR²R^{2a}-,
 -NR²C(O)CR²R^{2a}-, -CR²R^{2a}C(O)NR²-, -CR²R^{2a}NR²C(O)-, -NR²C(O)NR²-, -NR²-,
 -NR²CR²R^{2a}-, -CR²R^{2a}NR²-, O, -CR²R^{2a}O-, and -OCR²R^{2a}-;

Y is selected from one of the following rings and is substituted with 0-2 R^{4a};
 25 cyclopropyl, cyclopentyl, cyclohexyl, phenyl, piperidinyl, piperazinyl, pyridyl, pyrimidyl, furanyl, morpholinyl, thiophenyl, pyrrolyl, pyrrolidinyl, oxazolyl, isoxazolyl, isoxazolinyl, thiazolyl, isothiazolyl, pyrazolyl, imidazolyl, 1,2,3-oxadiazolyl, 1,2,4-oxadiazolyl, 1,2,5-oxadiazolyl, 1,3,4-oxadiazolyl, 1,2,3-thiadiazolyl, 1,2,4-thiadiazolyl, 1,2,5-thiadiazolyl, 1,3,4-thiadiazolyl, 1,2,3-triazolyl, 1,2,4-triazolyl, 1,2,5-triazolyl, 1,3,4-triazolyl, benzofuranyl,
 30 benzothiofuranyl, indolyl, benzimidazolyl, benzoxazolyl, benzthiazolyl, indazolyl, benzisoxazolyl, benzisothiazolyl, and isoindazolyl;

alternatively, Y is selected from the following bicyclic heteroaryl ring systems:



K is selected from O, S, NH, and N;

G_1 is absent or is selected from $-(CR^3R^{3a})_{1-3}-$, $-(CR^3R^{3a})_u C(O)(CR^3R^{3a})_{w-}$,

- 5 $-(CR^3R^{3a})_u O(CR^3R^{3a})_{w-}$, $-(CR^3R^{3a})_u NR^{3b}(CR^3R^{3a})_{w-}$,
 $-(CR^3R^{3a})_u C(O)NR^{3b}(CR^3R^{3a})_{w-}$, $-(CR^3R^{3a})_u NR^{3b}C(O)(CR^3R^{3a})_{w-}$,
 $-(CR^3R^{3a})_u NR^{3b}C(O)(CR^3R^{3a})_u C(O)NR^{3b}(CR^3R^{3a})_{w-}$, $-(CR^3R^{3a})_u S(CR^3R^{3a})_{w-}$,
 $-(CR^3R^{3a})_u S(O)(CR^3R^{3a})_{w-}$, $-(CR^3R^{3a})_u S(O)_2(CR^3R^{3a})_{w-}$,
 $-(CR^3R^{3a})_u S(O)NR^{3b}(CR^3R^{3a})_{w-}$, $-(CR^3R^{3a})_u NR^{3b}S(O)_2(CR^3R^{3a})_{w-}$, and
 10 $-(CR^3R^{3a})_u S(O)_2NR^{3b}(CR^3R^{3a})_{w-}$, wherein $u + w$ total 0, 1, or 2, provided that G_1
 does not form a N-S, NCH_2N , NCH_2O , or NCH_2S bond with either group to which it
 is attached;

- Z is selected from a bond, CH_2 , CH_2CH_2 , CH_2O , OCH_2 , $C(O)$, NH , CH_2NH ,
 $NHCH_2$, $CH_2C(O)$, $C(O)CH_2$, $C(O)NH$, $NHC(O)$, $NHC(O)CH_2C(O)NH$, $S(O)_2$,
 15 $CH_2S(O)_2$, $S(O)_2(CH_2)$, SO_2NH , and $NHSO_2$, provided that Z does not form a N-S,
 NCH_2N , NCH_2O , or NCH_2S bond with either group to which it is attached;

- R^{1a} , at each occurrence, is selected from H, $-(CH_2)_r-R^{1b}$, $-(CH(CH_3))_r-R^{1b}$,
 $-(C(CH_3)_2)_r-R^{1b}$, $-O-(CR^3R^{3a})_r-R^{1b}$, $-NR^2-(CR^3R^{3a})_r-R^{1b}$, and $-S-(CR^3R^{3a})_r-R^{1b}$,
 provided that R^{1a} forms other than an N-halo, N-S, O-O, or N-CN bond;
 20 alternatively, when two R^{1a} groups are attached to adjacent atoms, together
 with the atoms to which they are attached they form a 5-7 membered ring consisting
 of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and
 $S(O)_p$, this ring being substituted with 0-2 R^{4b} and 0-3 ring double bonds;

- R^{1b} is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, F, Cl, Br, I,
 25 CN, CHO, CF_3 , OR^2 , $-NR^2R^{2a}$, $-C(O)R^{2b}$, $-CO_2R^{2b}$, $-OC(O)R^2$, $-CO_2R^{2a}$,

-S(O)_pR^{2b}, -NR²(CH₂)_rOR², -NR²C(O)R^{2b}, -NR²C(O)NHR², -NR²C(O)₂R^{2a},
 -OC(O)NR²R^{2a}, -C(O)NR²R^{2a}, -C(O)NR²(CH₂)_rOR², -SO₂NR²R^{2a}, -NR²SO₂R²,
 C₅₋₆ carbocycle substituted with 0-2 R^{4b}, and 5-6 membered heterocycle substituted
 with 0-2 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from
 5 the group consisting of N, O, and S(O)_p, provided that R^{1b} forms other than an O-O,
 N-halo, N-S, or N-CN bond and provided that S(O)_pR² forms other than S(O)₂H or
 S(O)H;

R², at each occurrence, is selected from H, CF₃, CH₃, CH₂CH₃, CH₂CH₂CH₃,
 CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂, CH(CH₃)CH₂CH₃, C(CH₃)₃, benzyl,
 10 C₃₋₆ carbocycle substituted with 0-2 R^{4b}, C₃₋₆ carbocycle-CH₂- substituted with 0-2
 R^{4b}, and 5-6 membered heterocycle substituted with 0-2 R^{4b} and consisting of:
 carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and
 S(O)_p;

R^{2a}, at each occurrence, is selected from H, CF₃, CH₃, CH₂CH₃,
 15 CH₂CH₂CH₃, CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂, CH(CH₃)CH₂CH₃,
 C(CH₃)₃, benzyl, C₅₋₆ carbocycle substituted with 0-2 R^{4b}, and 5-6 membered
 heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4
 heteroatoms selected from the group consisting of N, O, and S(O)_p;

alternatively, R² and R^{2a}, together with the nitrogen atom to which they are
 20 attached, combine to form a 5 or 6 membered saturated, partially saturated, or
 unsaturated ring substituted with 0-2 R^{4b} and consisting of: 0-1 additional
 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{2b}, at each occurrence, is selected from CF₃, C₁₋₄ alkoxy, CH₃, CH₂CH₃,
 CH₂CH₂CH₃, CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂, CH(CH₃)CH₂CH₃,
 25 C(CH₃)₃, benzyl, C₅₋₆ carbocycle substituted with 0-2 R^{4b}, and 5-6 membered
 heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4
 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{2c}, at each occurrence, is selected from CF₃, OH, C₁₋₄ alkoxy, CH₃,
 CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂,
 30 CH(CH₃)CH₂CH₃, C(CH₃)₃, benzyl, C₅₋₆ carbocycle substituted with 0-2 R^{4b}, and

5-6 membered heterocycle substituted with 0-2 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^3 , at each occurrence, is selected from H, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, benzyl, and phenyl;

5 R^{3a} , at each occurrence, is selected from H, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, benzyl, and phenyl;

R^{3c} , at each occurrence, is selected from CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, benzyl, and phenyl;

R^{3d} , at each occurrence, is selected from H, CH₃, CH₂CH₃, CH₂CH₂CH₃,
10 CH(CH₃)₂, CH₂-phenyl, CH₂CH₂-phenyl, and C(=O) R^{3c} ;

R^4 , at each occurrence, is selected from H, =O, OR², -CH₂OR²,
-(CH₂)₂OR², F, Cl, Br, I, C₁₋₄ alkyl, CN, NO₂, -NR²R^{2a}, -CH₂NR²R^{2a},
-(CH₂)₂NR²R^{2a}, -C(O)R^{2c}, -NR²C(O)R^{2b}, -C(O)NR²R^{2a}, -SO₂NR²R^{2a}, -S(O)_pR^{5a},
CF₃, CF₂CF₃, 5-6 membered carbocycle substituted with 0-1 R^5 , and a 5-6 membered
15 heterocycle substituted with 0-1 R^5 and consisting of: carbon atoms and 1-4
heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{4a} , at each occurrence, is selected from H, =O, -(CR³R^{3a})_rOR²,
-(CR³R^{3a})_r-F, -(CR³R^{3a})_r-Br, -(CR³R^{3a})_r-Cl, C₁₋₄ alkyl, -(CR³R^{3a})_r-CN,
-(CR³R^{3a})_rNO₂, -(CR³R^{3a})_rNR²R^{2a}, -(CR³R^{3a})_rC(O)R^{2c}, -(CR³R^{3a})_rNR²C(O)R^{2b},
20 -(CR³R^{3a})_rC(O)NR²R^{2a}, -(CR³R^{3a})_rSO₂NR²R^{2a}, -(CR³R^{3a})_rNR²SO₂NR²R^{2a},
-(CR³R^{3a})_rNR²SO₂-C₁₋₄ alkyl, -(CR³R^{3a})_rC(O)NHSO₂-C₁₋₄ alkyl,
-(CR³R^{3a})_rNR²SO₂R⁵, -(CR³R^{3a})_rS(O)_pR⁵, -(CR³R^{3a})_r(CF₂)_rCF₃, phenyl substituted
with 0-1 R^5 , and a 5 membered aromatic heterocycle consisting of: carbon atoms and
1-3 heteroatoms selected from the group consisting of N, O, and S(O)_p substituted
25 with 0-1 R^5 ;

R^{4b} , at each occurrence, is selected from H, =O, OR³, CH₂OR³, F, Cl, CH₃,
CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂,
CH(CH₃)CH₂CH₃, C(CH₃)₃, CN, NO₂, -NR³R^{3a}, -CH₂NR³R^{3a}, -C(O)R³,
-CH₂-C(O)R³, -C(O)OR^{3c}, -CH₂C(O)OR^{3c}, -NR³C(O)R^{3a}, -CH₂NR³C(O)R^{3a},

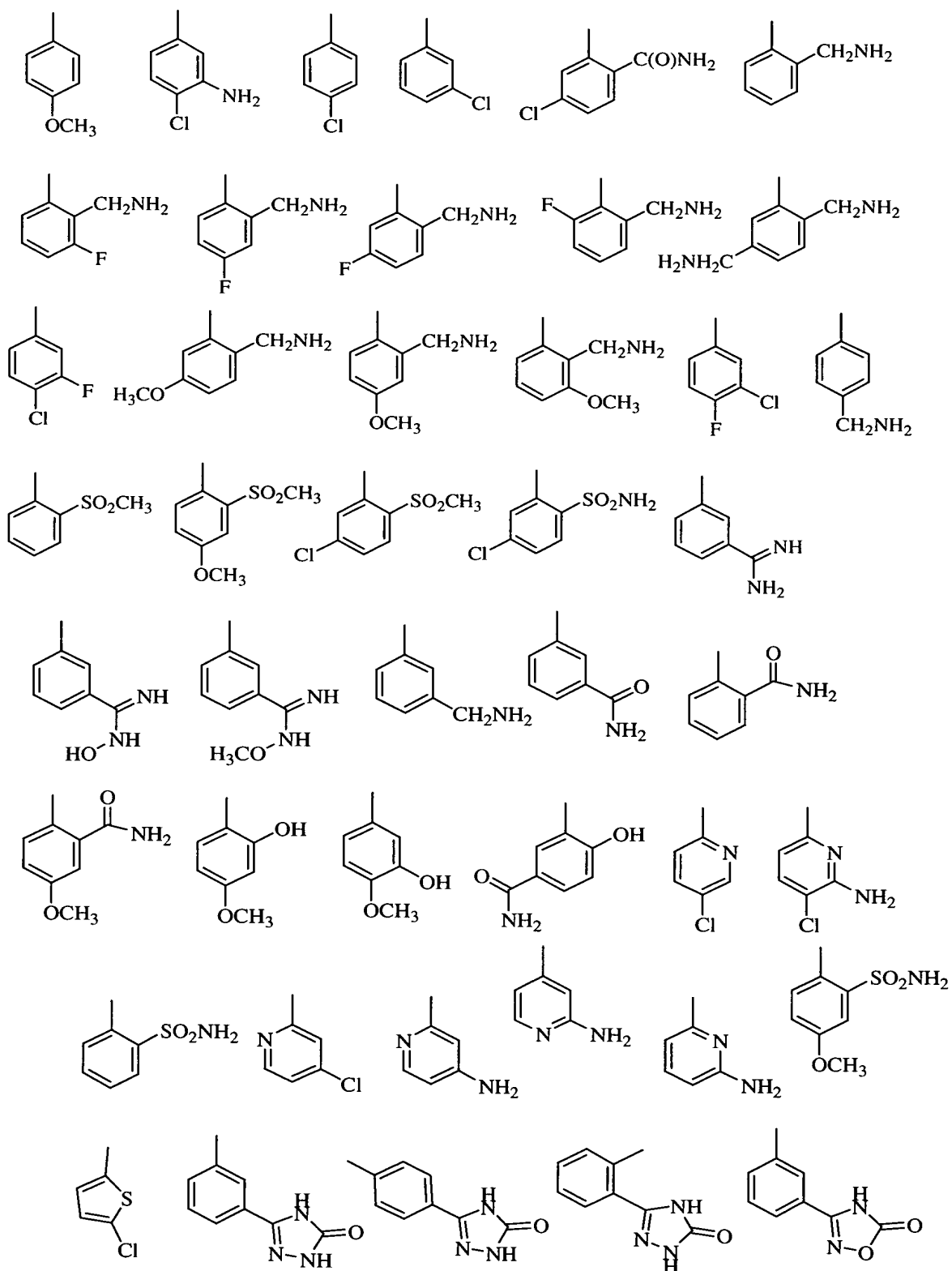
-C(O)NR³R^{3a}, -CH₂C(O)NR³R^{3a}, -NR³C(O)NR³R^{3a}, -CH₂NR³C(O)NR³R^{3a},
 -C(=NR³)NR³R^{3a}, -CH₂C(=NR³)NR³R^{3a}, -NR³C(=NR³)NR³R^{3a},
 -CH₂NR³C(=NR³)NR³R^{3a}, -SO₂NR³R^{3a}, -CH₂SO₂NR³R^{3a}, -NR³SO₂NR³R^{3a},
 -CH₂NR³SO₂NR³R^{3a}, -NR³SO₂-C₁₋₄ alkyl, -CH₂NR³SO₂-C₁₋₄ alkyl, -NR³SO₂CF₃,
 5 -CH₂NR³SO₂CF₃, -NR³SO₂-phenyl, -CH₂NR³SO₂-phenyl, -S(O)_pCF₃,
 -CH₂S(O)_pCF₃, -S(O)_p-C₁₋₄ alkyl, -CH₂S(O)_p-C₁₋₄ alkyl, -S(O)_p-phenyl,
 -CH₂S(O)_p-phenyl, CF₃, and CH₂-CF₃;

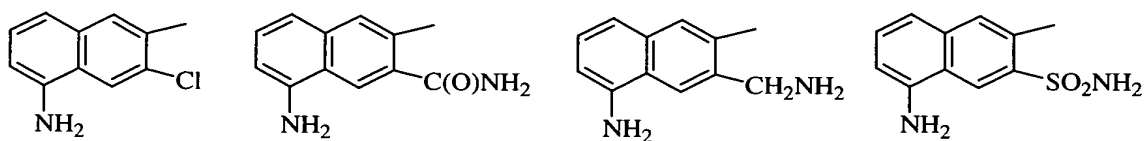
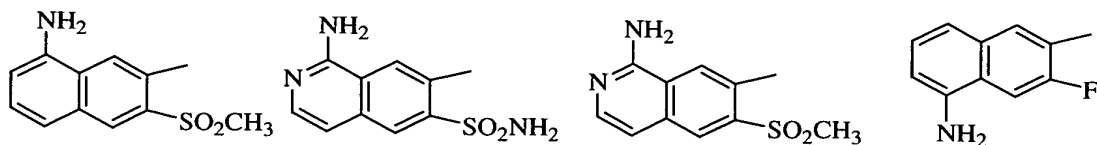
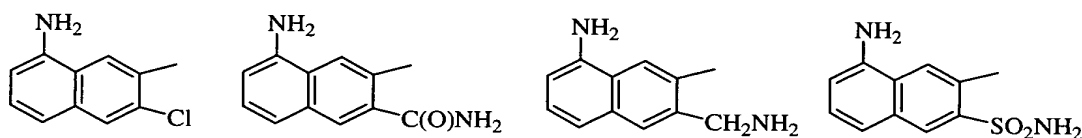
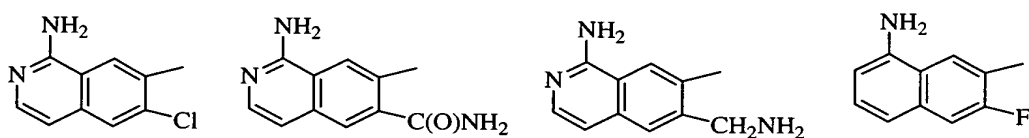
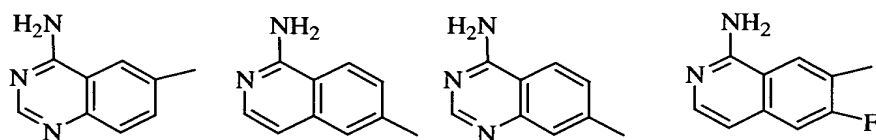
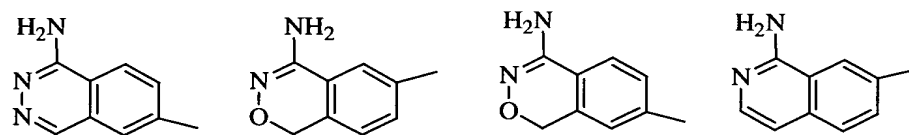
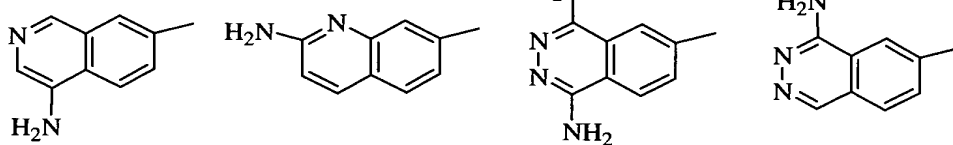
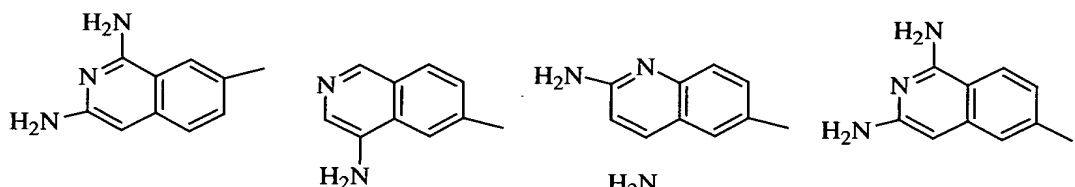
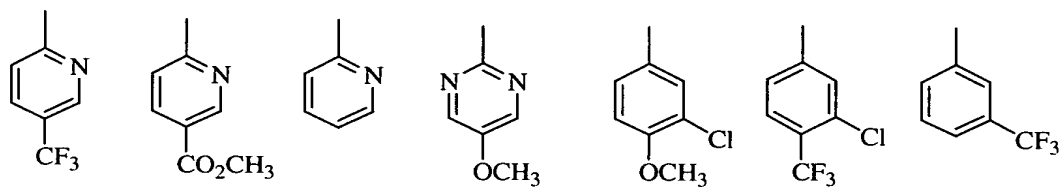
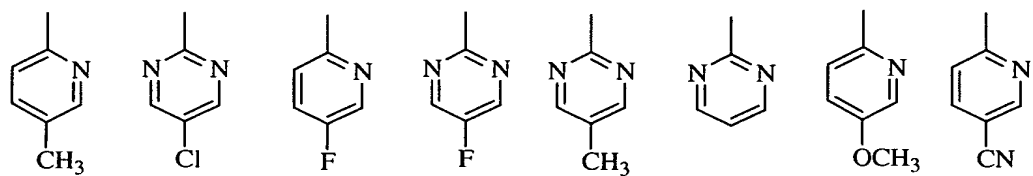
R⁵, at each occurrence, is selected from H, =O, CH₃, CH₂CH₃, CH₂CH₂CH₃,
 CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂, CH(CH₃)CH₂CH₃, C(CH₃)₃, OR³,
 10 -CH₂OR³, F, Cl, CN, NO₂, -NR³R^{3a}, -CH₂NR³R^{3a}, -C(O)R³, -CH₂C(O)R³,
 -C(O)OR^{3c}, -CH₂C(O)OR^{3c}, -NR³C(O)R^{3a}, -C(O)NR³R^{3a}, -NR³C(O)NR³R^{3a},
 -CH(=NOR^{3d}), -C(=NR³)NR³R^{3a}, -NR³C(=NR³)NR³R^{3a}, -SO₂NR³R^{3a},
 -NR³SO₂NR³R^{3a}, -NR³SO₂-C₁₋₄ alkyl, -NR³SO₂CF₃, -NR³SO₂-phenyl, -S(O)_pCF₃,
 -S(O)_p-C₁₋₄ alkyl, -S(O)_p-phenyl, CF₃, phenyl substituted with 0-2 R⁶, naphthyl
 15 substituted with 0-2 R⁶, and benzyl substituted with 0-2 R⁶;

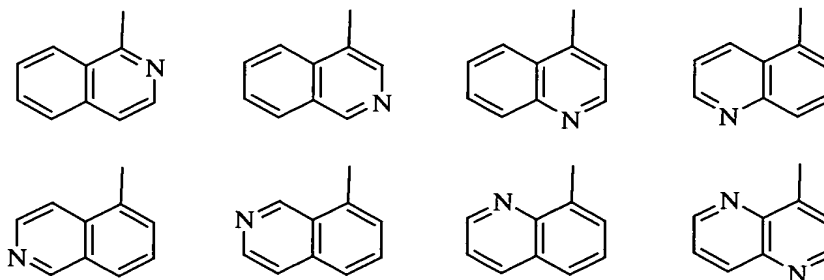
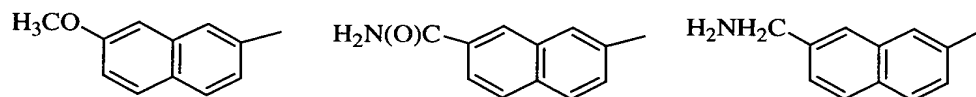
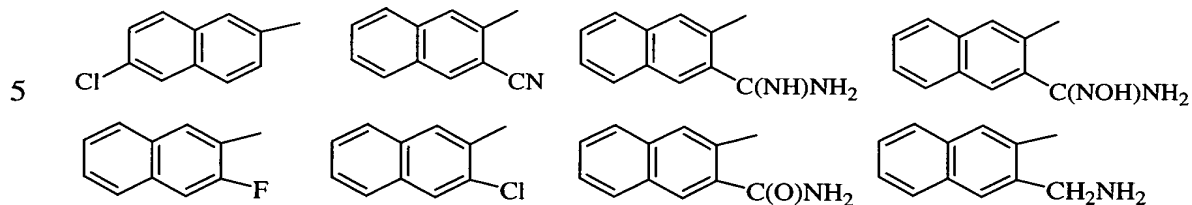
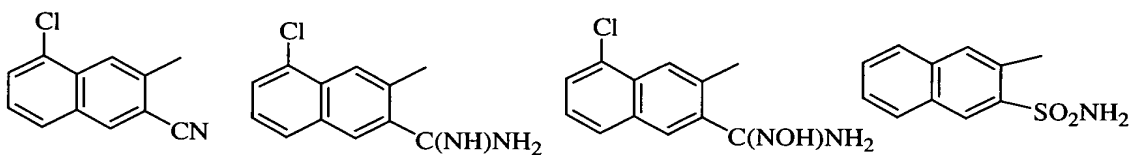
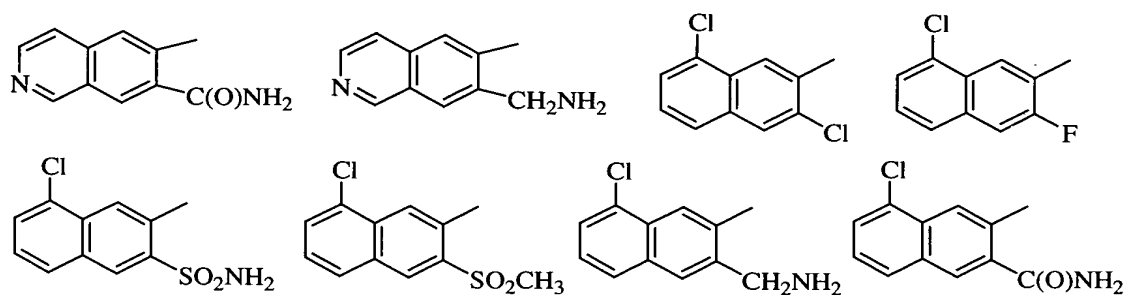
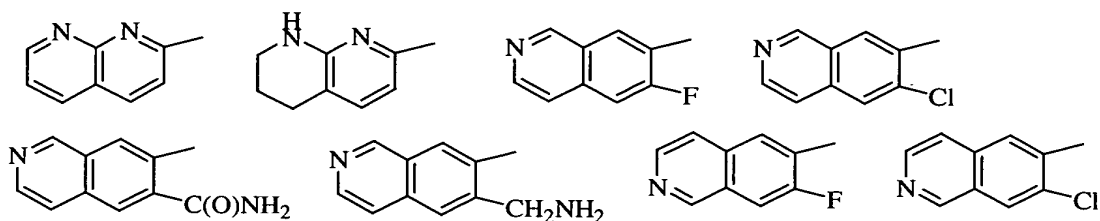
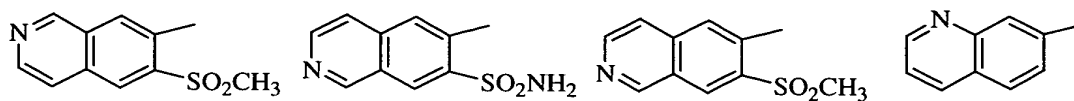
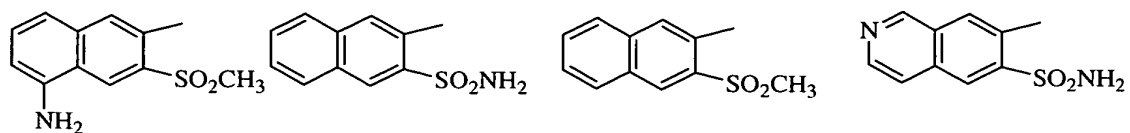
R⁶, at each occurrence, is selected from H, OH, OR², F, Cl, CH₃, CH₂CH₃,
 CH₂CH₂CH₃, CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂, CH(CH₃)CH₂CH₃,
 C(CH₃)₃, CN, NO₂, -NR²R^{2a}, -CH₂NR²R^{2a}, -C(O)R^{2b}, -CH₂C(O)R^{2b},
 -NR²C(O)R^{2b}, -NR²C(O)NR²R^{2a}, -C(=NH)NH₂, -NHC(=NH)NH₂, -SO₂NR²R^{2a},
 20 -NR²SO₂NR²R^{2a}, and -NR²SO₂C₁₋₄ alkyl; and

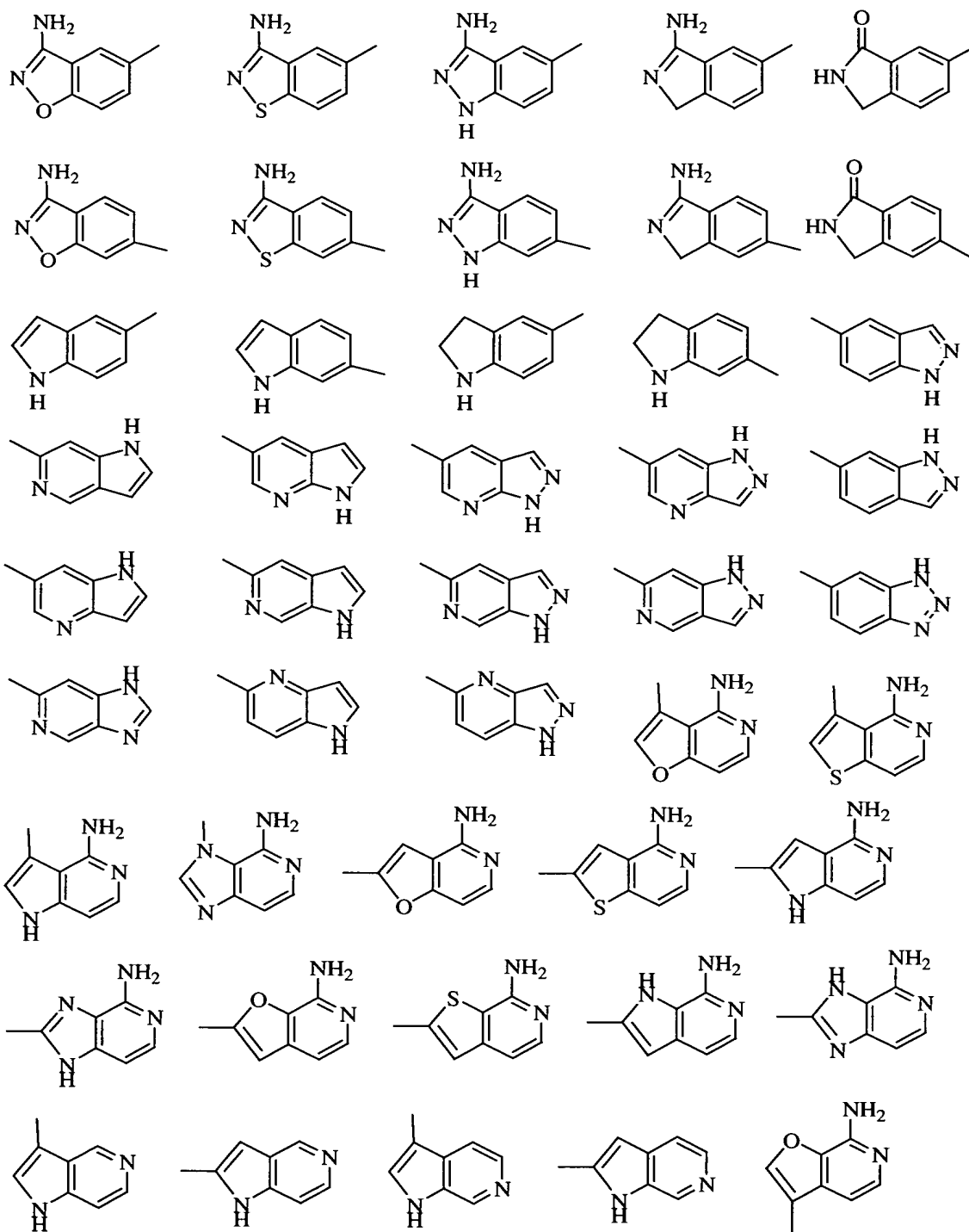
r, at each occurrence, is selected from 0, 1, 2, and 3.

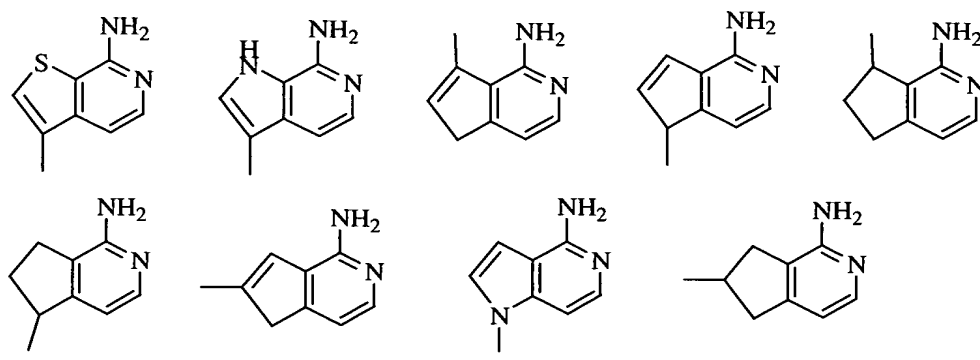
3. A compound according to Claim 2, wherein:
 25 ring M is selected from phenyl, pyrrole, furan, thiophene, pyrazole, imidazole,
 isoxazole, oxazole, isothiazole, thiazole, pyridine, and pyrimidine;
 ring M is substituted with 0-2 R^{1a};
 G is selected from the group:











R^{1a} , at each occurrence, is selected from H, R^{1b} , $-\text{CH}(\text{CH}_3)R^{1b}$, $-\text{C}(\text{CH}_3)_2R^{1b}$, $-\text{CH}_2R^{1b}$, and $-\text{CH}_2\text{CH}_2R^{1b}$, provided that R^{1a} forms other than an N-halo, N-S, or N-CN bond;

alternatively, when two R^{1a} groups are attached to adjacent atoms, together with the atoms to which they are attached, they form a 5-6 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and $\text{S}(\text{O})_p$, this ring being substituted with 0-2 R^{4b} and 0-3 ring double bonds;

R^{1b} is selected from H, CH_3 , CH_2CH_3 , F, Cl, Br, CN, CHO, CF_3 , OR^2 , $-\text{NR}^2R^{2a}$, $-\text{C}(\text{O})R^{2b}$, $-\text{CO}_2R^{2b}$, $-\text{OC}(\text{O})R^2$, $-\text{CO}_2R^{2a}$, $-\text{S}(\text{O})_pR^2$, $-\text{NR}^2(\text{CH}_2)_r\text{OR}^2$, $-\text{NR}^2\text{C}(\text{O})R^{2b}$, $-\text{C}(\text{O})\text{NR}^2R^{2a}$, $-\text{SO}_2\text{NR}^2R^{2a}$, $-\text{NR}^2\text{SO}_2R^2$, phenyl substituted with 0-2 R^{4b} , and 5-6 membered aromatic heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $\text{S}(\text{O})_p$ and substituted with 0-2 R^{4b} , provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond;

R^2 , at each occurrence, is selected from H, CF_3 , CH_3 , CH_2CH_3 , $\text{CH}_2\text{CH}_2\text{CH}_3$, $\text{CH}(\text{CH}_3)_2$, phenyl substituted with 0-2 R^{4b} , a benzyl substituted with 0-2 R^{4b} , C_{3-6} cycloalkyl substituted with 0-2 R^{4b} , C_{3-6} cycloalkyl- CH_2 - substituted with 0-2 R^{4b} , and 5-6 membered aromatic heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $\text{S}(\text{O})_p$;

R^{2a} , at each occurrence, is selected from H, CF_3 , CH_3 , CH_2CH_3 , $\text{CH}_2\text{CH}_2\text{CH}_3$, $\text{CH}(\text{CH}_3)_2$, benzyl, phenyl substituted with 0-2 R^{4b} , and 5-6 membered

aromatic heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{2b} , at each occurrence, is selected from CF₃, C₁₋₄ alkoxy, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, benzyl, phenyl substituted with 0-2 R^{4b} , and 5-6 membered

5 aromatic heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{2c} , at each occurrence, is selected from CF₃, OH, OCH₃, OCH₂CH₃, OCH₂CH₂CH₃, OCH(CH₃)₂, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, benzyl, phenyl substituted with 0-2 R^{4b} , and 5-6 membered aromatic heterocycle substituted with 0-2 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

alternatively, R^2 and R^{2a} , together with the nitrogen atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-2 R^{4b} and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^4 , at each occurrence, is selected from H, -(CH₂)₂OR², -CH₂OR², OR², F, Cl, Br, I, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, CH₂CH₂CH₂CH₃, CH₂CH(CH₃)₂, CH(CH₃)CH₂CH₃, C(CH₃)₃, CN, NO₂, -NR²R^{2a}, -CH₂NR²R^{2a}, -(CH₂)₂NR²R^{2a}, -C(O)R^{2c}, -NR²C(O)R^{2b}, -C(O)NR²R^{2a}, -SO₂NR²R^{2a}, CF₃, and CF₂CF₃;

R^{4a} , at each occurrence, is selected from H, =O, -(CH₂)_rOR², -(CH₂)_r-F, -(CH₂)_r-Br, -(CH₂)_r-Cl, C₁₋₄ alkyl, -(CH₂)_r-CN, -(CH₂)_rNO₂, -(CH₂)_rNR²R^{2a}, -(CH₂)_rC(O)R^{2c}, -(CH₂)_rNR²C(O)R^{2b}, -(CH₂)_rC(O)NR²R^{2a}, -(CH₂)_rSO₂NR²R^{2a}, -(CH₂)_rNR²SO₂NR²R^{2a}, -(CH₂)_rNR²SO₂-C₁₋₄ alkyl, -(CH₂)_rC(O)NHSO₂-C₁₋₄ alkyl, -(CH₂)_rNR²SO₂R⁵, -(CH₂)_rS(O)_pR⁵, -(CH₂)_r(CF₂)_rCF₃, phenyl substituted with 0-1 R⁵, and a 5 membered aromatic heterocycle consisting of: carbon atoms and 1-3 heteroatoms selected from the group consisting of N, O, and S(O)_p substituted with 0-1 R⁵;

R^{4b} , at each occurrence, is selected from H, =O, OR^3 , $-CH_2OR^3$, F, Cl, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, CN, NO_2 , $-NR^3R^{3a}$, $-CH_2NR^3R^{3a}$, $-C(O)R^3$, $-CH_2-C(O)R^3$, $-C(O)OR^{3c}$, $-CH_2-C(O)OR^{3c}$, $-NR^3C(O)R^{3a}$, $-CH_2NR^3C(O)R^{3a}$, $-C(O)NR^3R^{3a}$, $-CH_2-C(O)NR^3R^{3a}$, $-SO_2NR^3R^{3a}$, $-CH_2SO_2NR^3R^{3a}$,

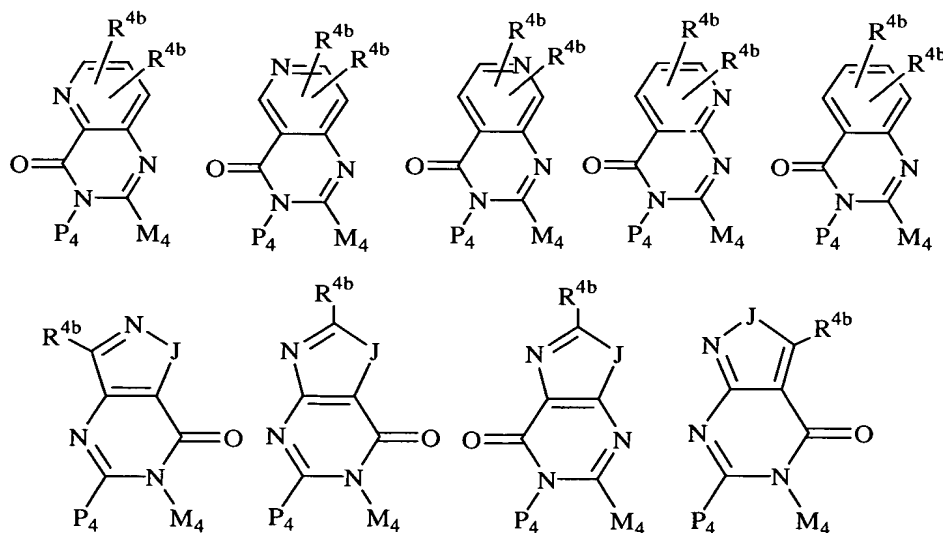
- 5 $-NR^3SO_2-C_{1-4}$ alkyl, $-CH_2NR^3SO_2-C_{1-4}$ alkyl, $-NR^3SO_2$ -phenyl, $-CH_2NR^3SO_2$ -phenyl, $-S(O)_pCF_3$, $-CH_2S(O)_pCF_3$, $-S(O)_p-C_{1-4}$ alkyl, $-CH_2S(O)_p-C_{1-4}$ alkyl, $-S(O)_p$ -phenyl, $-CH_2S(O)_p$ -phenyl, and CF_3 ;

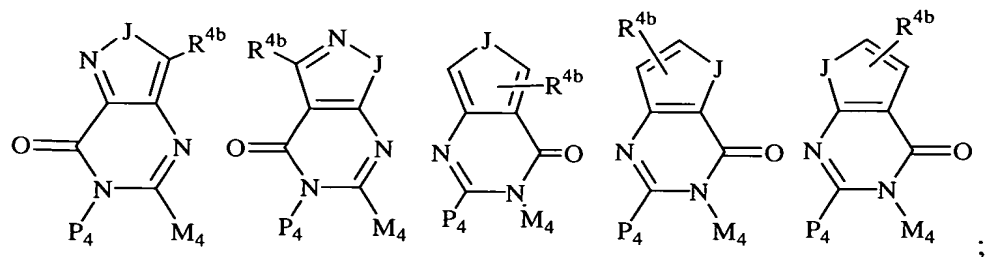
- R^5 , at each occurrence, is selected from H, =O, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, OR^3 , CH_2OR^3 , F, Cl, CN, NO_2 , $-NR^3R^{3a}$, $-CH_2NR^3R^{3a}$, $-C(O)R^3$, $-CH_2C(O)R^3$, $-C(O)OR^{3c}$, $-CH_2C(O)OR^{3c}$, $-NR^3C(O)R^{3a}$, $-C(O)NR^3R^{3a}$, $-SO_2NR^3R^{3a}$, $-NR^3SO_2-C_{1-4}$ alkyl, $-NR^3SO_2CF_3$, $-NR^3SO_2$ -phenyl, $-S(O)_pCF_3$, $-S(O)_p-C_{1-4}$ alkyl, $-S(O)_p$ -phenyl, CF_3 , phenyl substituted with 0-2 R^6 , naphthyl substituted with 0-2 R^6 , and benzyl substituted with 0-2 R^6 ;

- R^6 , at each occurrence, is selected from H, OH, OR^2 , F, Cl, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, CN, NO_2 , $-NR^2R^{2a}$, $-CH_2NR^2R^{2a}$, $-C(O)R^{2b}$, $-CH_2C(O)R^{2b}$, $-NR^2C(O)R^{2b}$, $-SO_2NR^2R^{2a}$, and $NR^2SO_2C_{1-4}$ alkyl; and

r, at each occurrence, is selected from 0, 1, and 2.

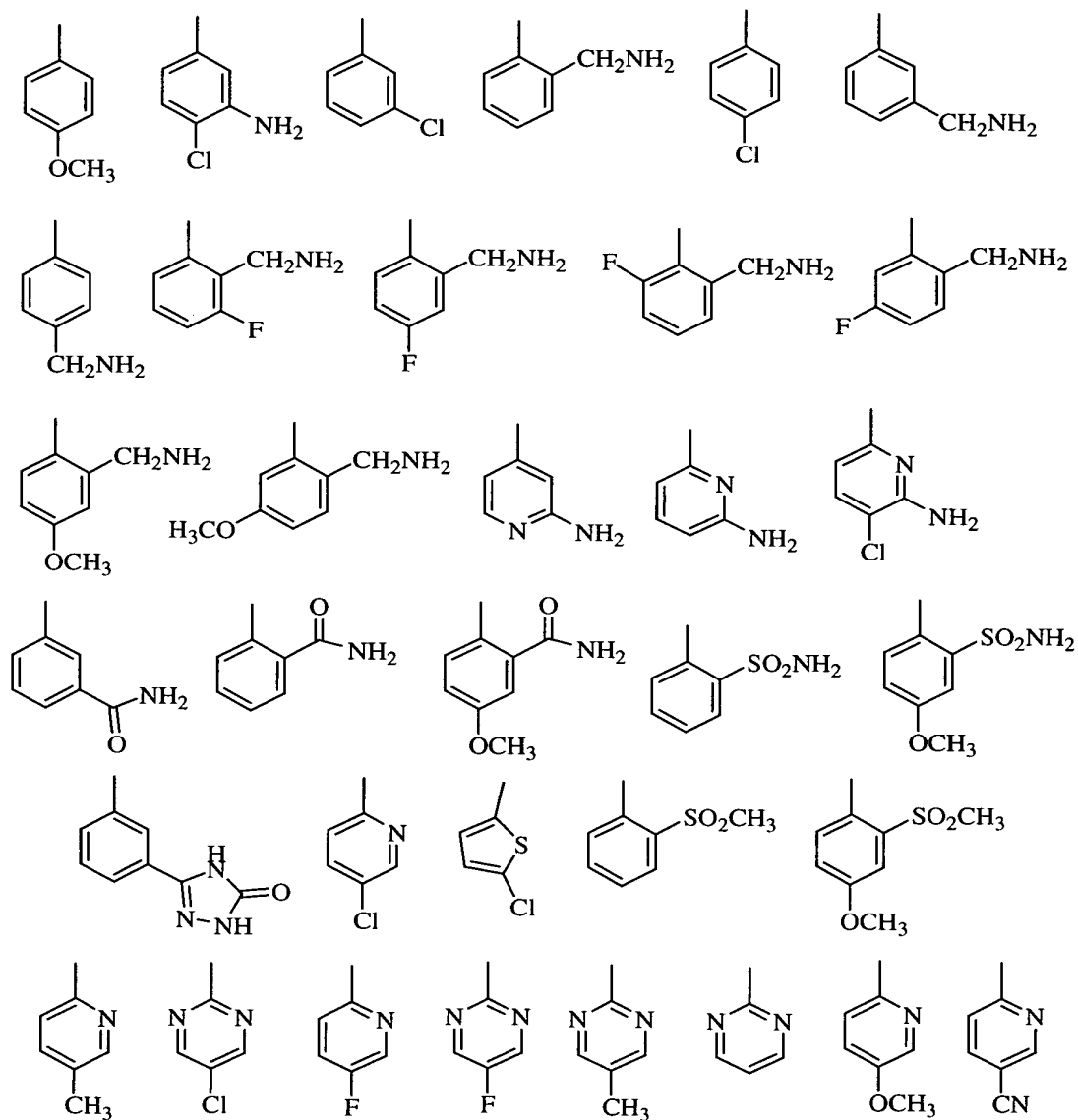
- 20 4. A compound according to Claim 3, wherein the compound is selected from:



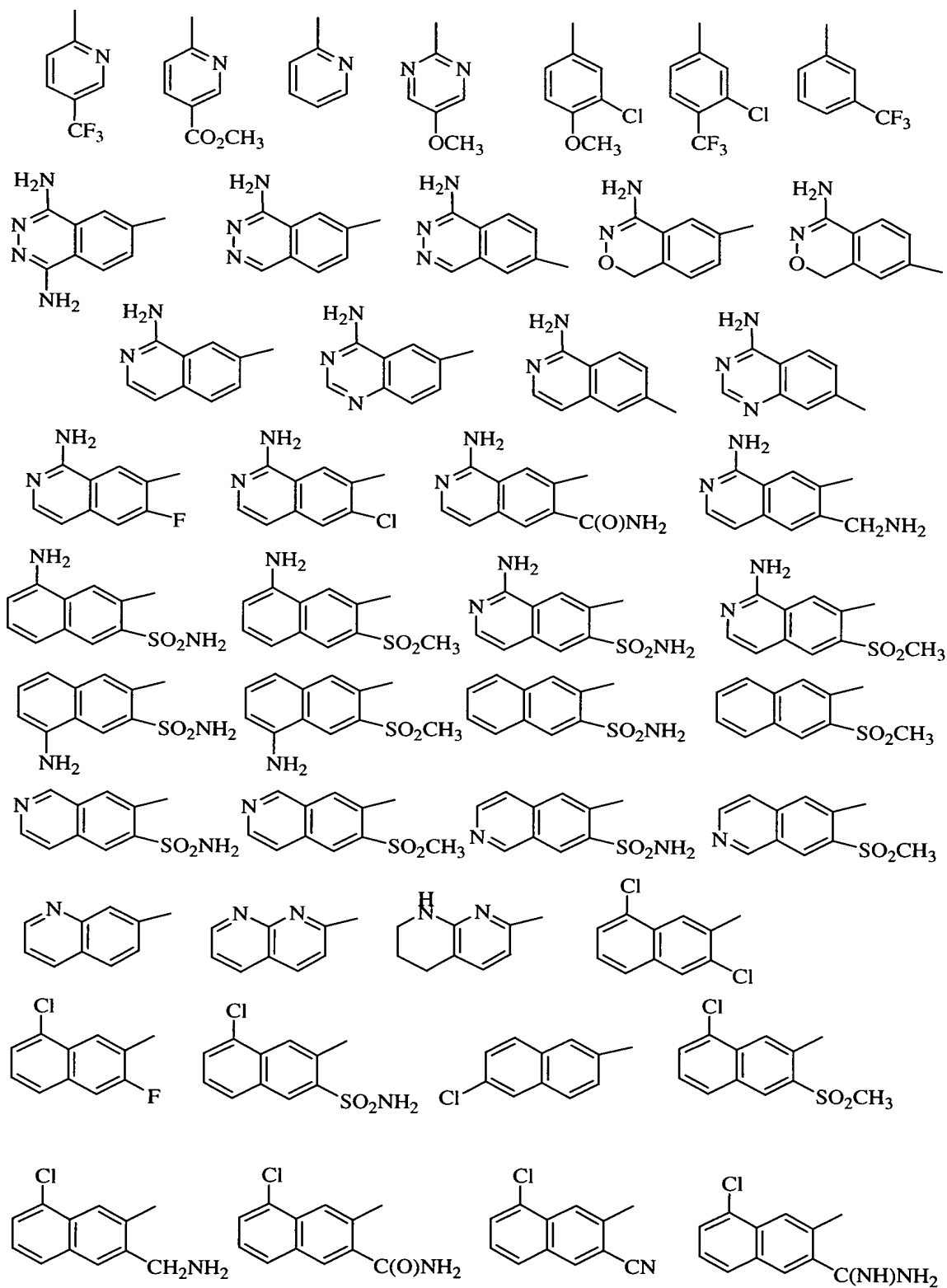


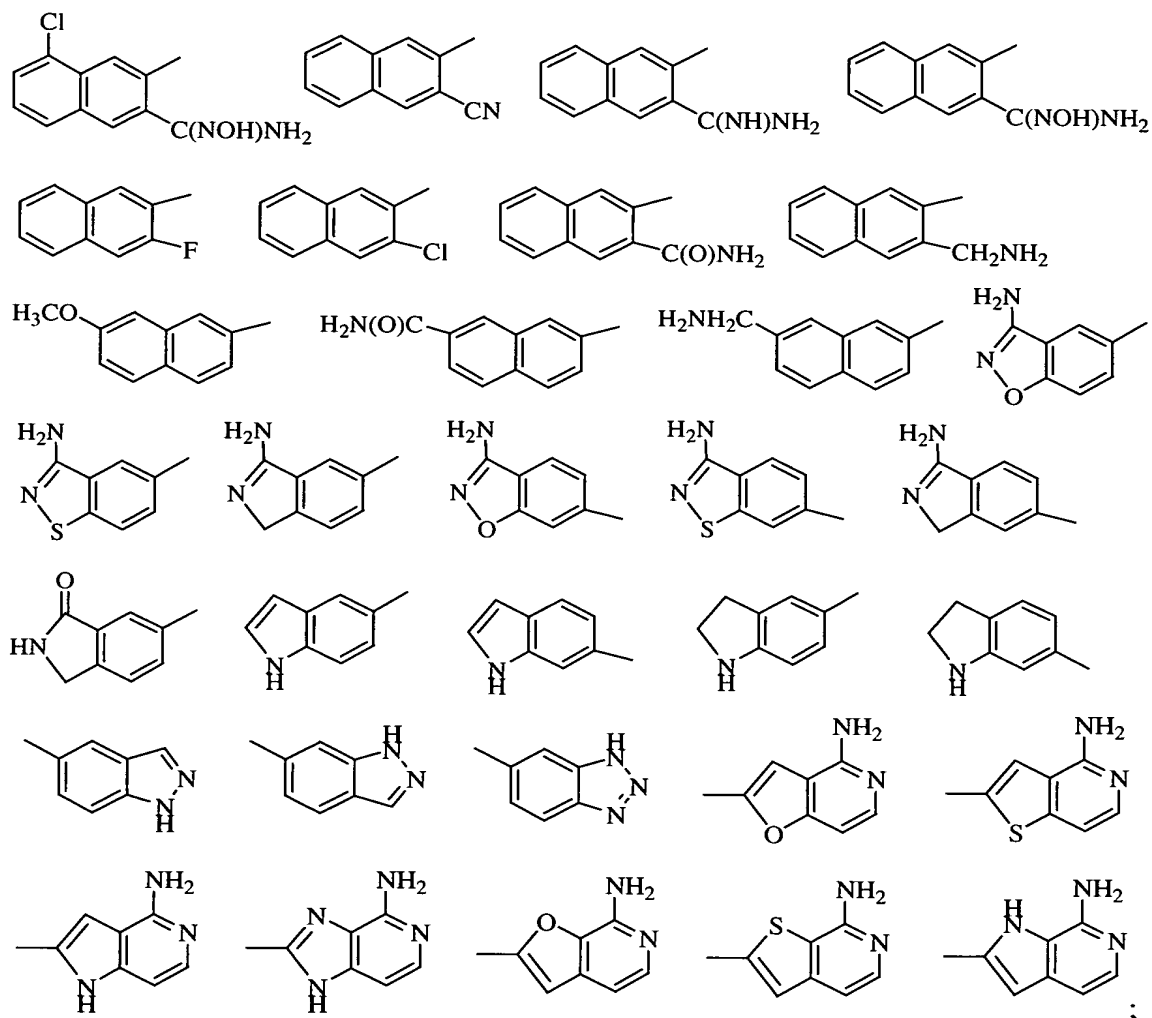
J is selected from O, S, NH, and NR^{1a};

G is selected from the group:



5





5 R^{1a}, at each occurrence, is selected from H, R^{1b}, -CH(CH₃)R^{1b}, -C(CH₃)₂R^{1b}, and -CH₂R^{1b}, provided that R^{1a} forms other than an N-halo, N-S, or N-CN bond;

R^{1b} is selected from CH_3 , CH_2CH_3 , F, Cl, Br, -CN, CF_3 , OR^2 , $-NR^2R^{2a}$, $-C(O)R^{2b}$, $-CO_2R^{2b}$, $-CO_2R^{2a}$, $-S(O)_pR^2$, $-C(O)NR^2R^{2a}$, $-SO_2NR^2R^{2a}$, $-NR^2SO_2R^2$, and 5-6 membered aromatic heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-2 R^{4b} , provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond;

R², at each occurrence, is selected from H, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, phenyl substituted with 0-1 R^{4b}, benzyl substituted with 0-1 R^{4b},

C₃₋₅ cycloalkyl substituted with 0-1 R^{4b}, C₃₋₅ cycloalkyl-CH₂- substituted with 0-1 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-1 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

5 R^{2a}, at each occurrence, is selected from H, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, benzyl, phenyl substituted with 0-1 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-1 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

alternatively, R² and R^{2a}, together with the nitrogen atom to which they are
10 attached, combine to form a 5 or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-1 R^{4b} and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{2b}, at each occurrence, is selected from OCH₃, -OCH₂CH₃, OCH₂CH₂CH₃, OCH(CH₃)₂, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, benzyl, phenyl substituted
15 with 0-1 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-1 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{2c}, at each occurrence, is selected from OH, OCH₃, OCH₂CH₃, OCH₂CH₂CH₃, OCH(CH₃)₂, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, benzyl,
20 phenyl substituted with 0-1 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-1 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

R⁴, at each occurrence, is selected from OH, OR², CH₂OR², (CH₂)₂OR², F, Br, Cl, I, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, CH₂CH₂CH₂CH₃,
25 CH₂CH(CH₃)₂, CH(CH₃)CH₂CH₃, C(CH₃)₃, -NR²R^{2a}, -CH₂NR²R^{2a}, -(CH₂)₂NR²R^{2a}, CF₃, and CF₂CF₃;

R^{4a}, at each occurrence, is selected from H, =O, -(CH₂)_rOR², F, Br, Cl, C₁₋₄ alkyl, -(CH₂)_rNR²R^{2a}, -(CH₂)_rC(O)R^{2c}, -(CH₂)_rNR²C(O)R^{2b},
-(CH₂)_rC(O)NR²R^{2a}, -(CH₂)_rSO₂NR²R^{2a}, -(CH₂)_rNR²SO₂R⁵, -(CH₂)_rS(O)_pR⁵,

$-(\text{CH}_2)_r(\text{CF}_2)_r\text{CF}_3$, phenyl substituted with 0-1 R^5 , and a 5 membered aromatic heterocycle consisting of: carbon atoms and 1-3 N and is substituted with 1 R^5 ;

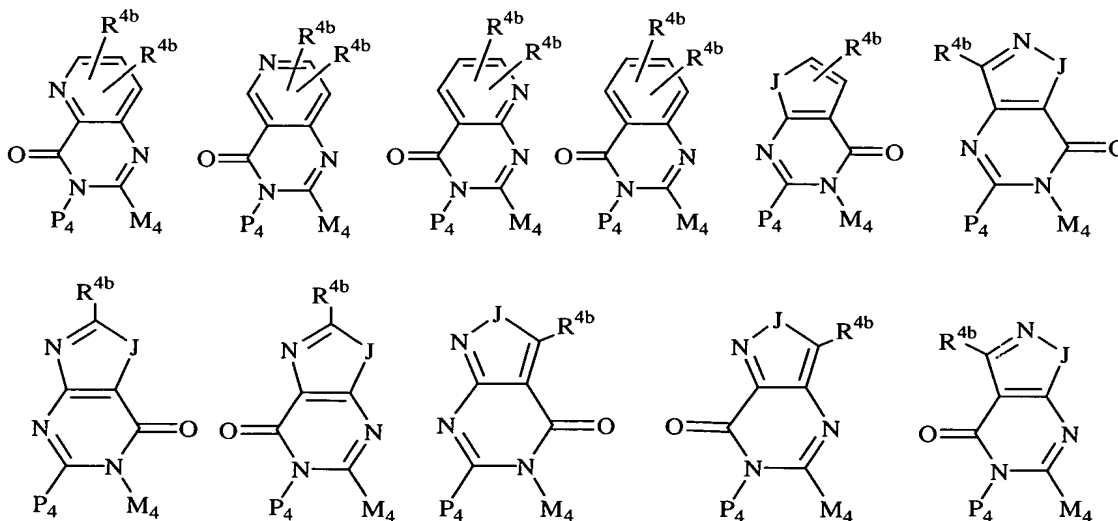
R^{4b} , at each occurrence, is selected from H, =O, OR^3 , $-\text{CH}_2\text{OR}^3$, F, Cl, CH_3 , CH_2CH_3 , $\text{CH}_2\text{CH}_2\text{CH}_3$, $\text{CH}(\text{CH}_3)_2$, CN, NO_2 , $-\text{NR}^3\text{R}^{3a}$, $-\text{CH}_2\text{NR}^3\text{R}^{3a}$, $-\text{C}(\text{O})\text{R}^3$,
 5 $-\text{C}(\text{O})\text{OR}^{3c}$, $-\text{NR}^3\text{C}(\text{O})\text{R}^{3a}$, $-\text{C}(\text{O})\text{NR}^3\text{R}^{3a}$, $-\text{SO}_2\text{NR}^3\text{R}^{3a}$, $-\text{NR}^3\text{SO}_2-\text{C}_{1-4}$ alkyl, $-\text{NR}^3\text{SO}_2$ -phenyl, $-\text{S}(\text{O})_p-\text{C}_{1-4}$ alkyl, $-\text{S}(\text{O})_p$ -phenyl, and CF_3 ;

R^5 , at each occurrence, is selected from H, =O, CH_3 , CH_2CH_3 , $\text{CH}_2\text{CH}_2\text{CH}_3$, $\text{CH}(\text{CH}_3)_2$, OR^3 , CH_2OR^3 , F, Cl, CN, NO_2 , $-\text{NR}^3\text{R}^{3a}$, $-\text{CH}_2\text{NR}^3\text{R}^{3a}$, $-\text{C}(\text{O})\text{R}^3$, $-\text{C}(\text{O})\text{OR}^{3c}$, $-\text{NR}^3\text{C}(\text{O})\text{R}^{3a}$, $-\text{C}(\text{O})\text{NR}^3\text{R}^{3a}$, $-\text{SO}_2\text{NR}^3\text{R}^{3a}$, $-\text{NR}^3\text{SO}_2-\text{C}_{1-4}$ alkyl,
 10 $-\text{NR}^3\text{SO}_2$ -phenyl, $-\text{S}(\text{O})_p-\text{C}_{1-4}$ alkyl, $-\text{S}(\text{O})_p$ -phenyl, CF_3 , phenyl substituted with 0-2 R^6 , naphthyl substituted with 0-2 R^6 , and benzyl substituted with 0-2 R^6 ; and

R^6 , at each occurrence, is selected from H, OH, OR^2 , F, Cl, CH_3 , CH_2CH_3 , $\text{CH}_2\text{CH}_2\text{CH}_3$, $\text{CH}(\text{CH}_3)_2$, CN, NO_2 , $-\text{NR}^2\text{R}^{2a}$, $-\text{CH}_2\text{NR}^2\text{R}^{2a}$, $-\text{C}(\text{O})\text{R}^{2b}$, $-\text{CH}_2\text{C}(\text{O})\text{R}^{2b}$, $-\text{NR}^2\text{C}(\text{O})\text{R}^{2b}$, and $-\text{SO}_2\text{NR}^2\text{R}^{2a}$.

15

5. A compound according to Claim 4, wherein the compound is selected from:



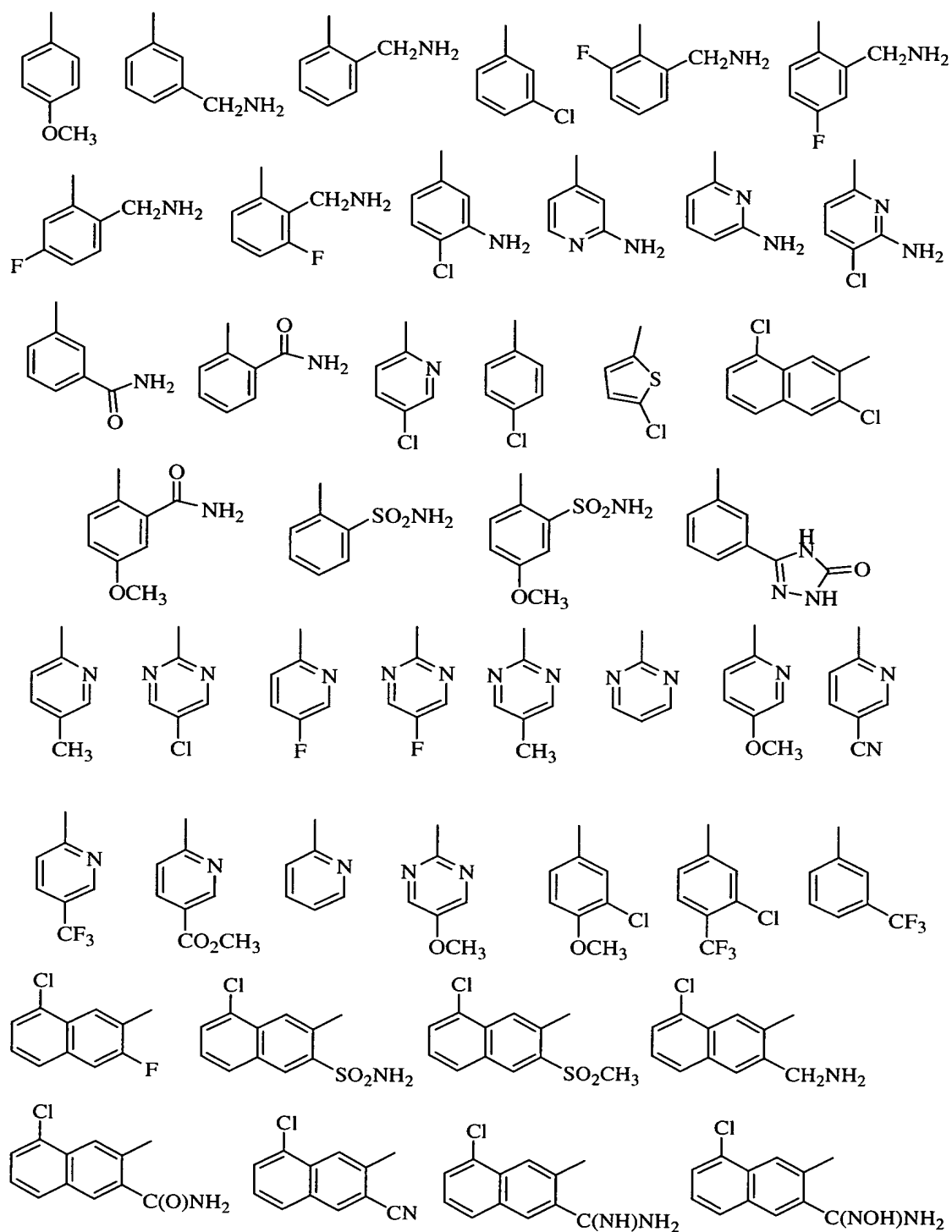
J is selected from O, S, NH, and NR^{1a} ;

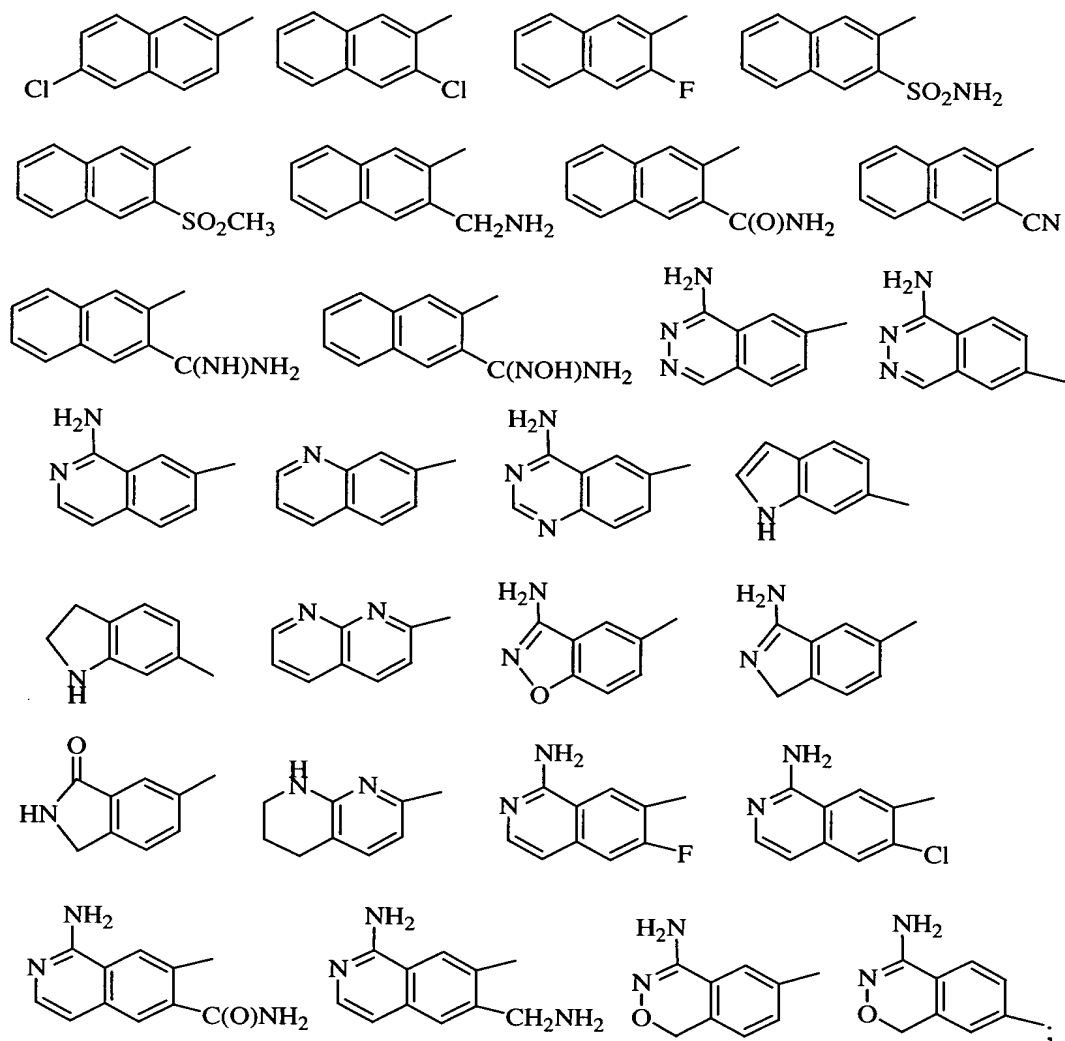
P_4 is $-\text{G}_1-\text{G}$;

20

M_4 is $-\text{Z}-\text{A}-\text{B}$;

G is selected from:





G_1 is absent or is selected from CH_2CH_2 , CH_2O , OCH_2 , CH_2NH , $NHCH_2$, $CH_2C(O)$, $C(O)CH_2$, $C(O)NH$, $NHC(O)$, SO_2NH , and $NHSO_2$, provided that G_1 does not form a N-S, NCH_2N , NCH_2O , or NCH_2S bond with either group to which it is attached;

A is selected from phenyl, pyridyl, and pyrimidyl, and is substituted with 0-2 R^4 ;

B is selected from phenyl, pyrrolidinyl, N-pyrrolidino-carbonyl, morpholinyl, N-morpholino-carbonyl, 1,2,3-triazolyl, imidazolyl, and benzimidazolyl, and is substituted with 0-1 R^{4a} ;

R^{1a} , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, CH_2F , CH_2Cl , Br, CH_2Br , $-CN$, CH_2CN , CF_3 , CH_2CF_3 , OCH_3 , CH_2OH ,

$C(CH_3)_2OH$, CH_2OCH_3 , NH_2 , CH_2NH_2 , $NHCH_3$, CH_2NHCH_3 , $N(CH_3)_2$,
 $CH_2N(CH_3)_2$, CO_2H , $COCH_3$, CO_2CH_3 , $CH_2CO_2CH_3$, SCH_3 , CH_2SCH_3 , $S(O)CH_3$,
 $CH_2S(O)CH_3$, $S(O)_2CH_3$, $CH_2S(O)_2CH_3$, $C(O)NH_2$, $CH_2C(O)NH_2$, SO_2NH_2 ,
 $CH_2SO_2NH_2$, $NHSO_2CH_3$, $CH_2NHSO_2CH_3$, pyridin-2-yl, pyridin-3-yl, pyridin-4-yl,
 5 pyridin-2-yl-N-oxide, pyridin-3-yl-N-oxide, pyridin-4-yl-N-oxide, imidazol-1-yl,
 CH_2 -imidazol-1-yl, 4-methyl-oxazol-2-yl, 4-N,N-dimethylaminomethyl-oxazol-2-yl,
 1,2,3,4-tetrazol-1-yl, 1,2,3,4-tetrazol-5-yl, $-CH_2$ -1,2,3,4-tetrazol-1-yl, and
 $-CH_2$ -1,2,3,4-tetrazol-5-yl, provided that R^{1a} forms other than an N-halo, N-S, or
 N-CN bond;

10 R^2 , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , cyclopropylmethyl,
 cyclobutyl, and cyclopentyl;

R^{2a} , at each occurrence, is H or CH_3 ;

alternatively, R^2 and R^{2a} , together with the atom to which they are attached,
 combine to form pyrrolidine substituted with 0-2 R^{4b} or piperidine substituted with
 15 0-2 R^{4b} ;

R^{2b} , at each occurrence, is selected from OCH_3 , OCH_2CH_3 , CH_3 , and
 CH_2CH_3 ;

R^{2c} , at each occurrence, is selected from OH, OCH_3 , OCH_2CH_3 , CH_3 , and
 CH_2CH_3 ;

20 R^4 , at each occurrence, is selected from OH, OR^2 , CH_2OR^2 , $(CH_2)_2OR^2$, F,
 Br, Cl, I, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$,
 $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, $-NR^2R^{2a}$, $-CH_2NR^2R^{2a}$,
 $-(CH_2)_2NR^2R^{2a}$, CF_3 , and CF_2CF_3 ;

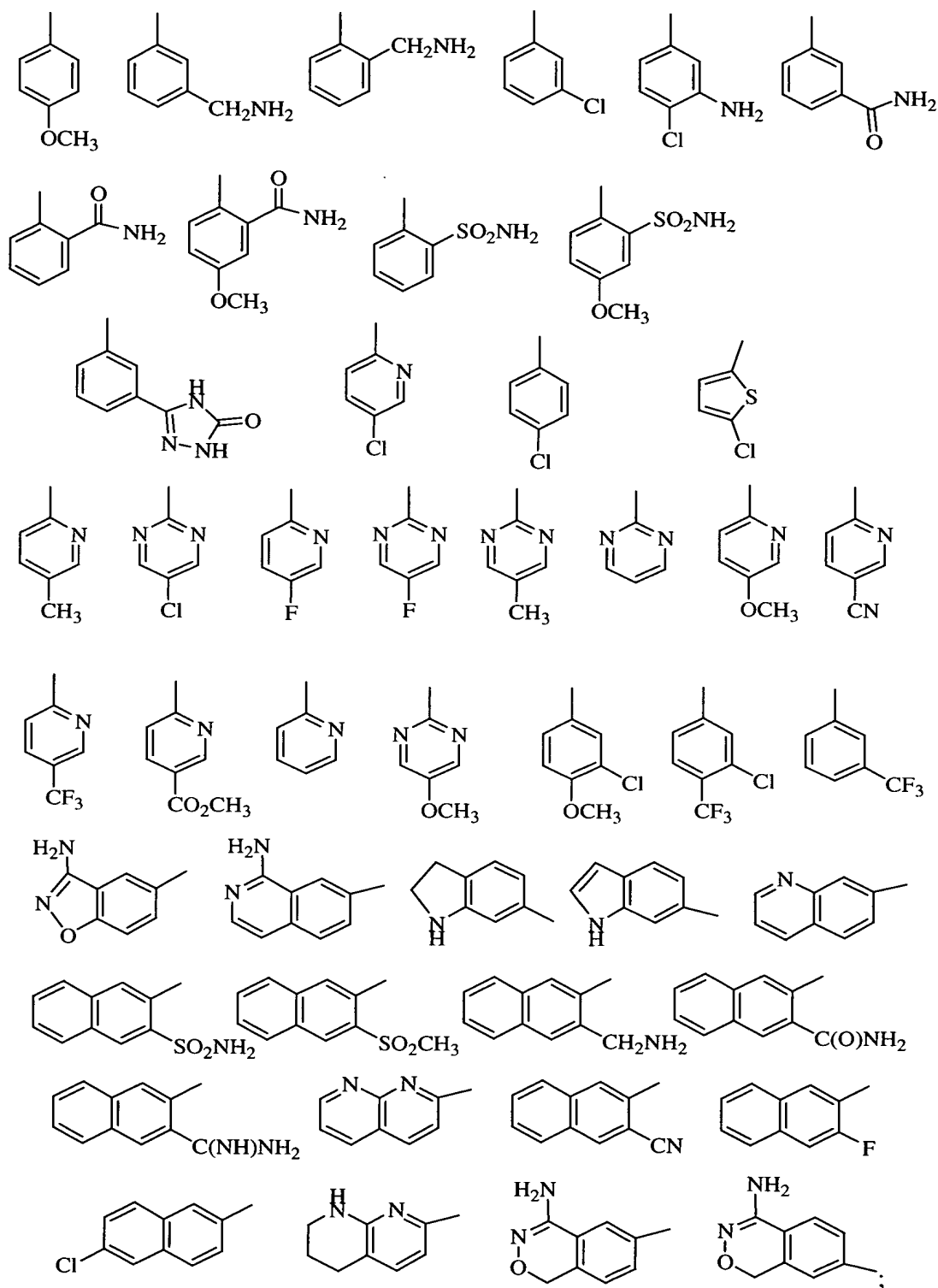
R^{4a} is selected from C_{1-4} alkyl, CF_3 , OR^2 , $-CH_2OR^2$, $-(CH_2)_2OR^2$, $-NR^2R^{2a}$,
 25 $-CH_2NR^2R^{2a}$, $-(CH_2)_2NR^2R^{2a}$, $-S(O)_pR^5$, $-SO_2NR^2R^{2a}$, and 1- CF_3 -tetrazol-2-yl;

R^{4b} , at each occurrence, is selected from H, CH_3 , and OH; and

R^5 , at each occurrence, is selected from CF_3 , CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$,
 $CH(CH_3)_2$, phenyl, and benzyl.

6. A compound according to Claim 5, wherein the compound is selected from:

G is selected from:



A is selected from the group: phenyl, 2-pyridyl, 3-pyridyl, 2-pyrimidyl, 2-Cl-phenyl, 3-Cl-phenyl, 2-F-phenyl, 3-F-phenyl, 2-methylphenyl, 2-aminophenyl, and 2-methoxyphenyl; and

B is selected from the group: 2-(aminosulfonyl)phenyl,
 5 2-(methylaminosulfonyl)phenyl, N-pyrrolidino-carbonyl, 2-(methylsulfonyl)phenyl, 2-(N,N-dimethylaminomethyl)phenyl, 2-(N-methylaminomethyl)phenyl, 2-(N-ethyl-N-methylaminomethyl)phenyl, 2-(N-pyrrolidinylmethyl)phenyl, 1-methyl-2-imidazolyl, 2-methyl-1-imidazolyl, 2-(dimethylaminomethyl)-1-imidazolyl, 2-(methylaminomethyl)-1-imidazolyl,
 10 2-(N-(cyclopropylmethyl)aminomethyl)phenyl, 2-(N-(cyclobutyl)aminomethyl)phenyl, 2-(N-(cyclopentyl)aminomethyl)phenyl, 2-(N-(4-hydroxypiperidinyl)methyl)phenyl, and 2-(N-(3-hydroxypyrrolidinyl)methyl)phenyl.

15

7. A compound according to Claim 6, wherein the compound is selected from the group:

2-biphenyl-4-yl-6-chloro-3-(4-chloro-phenyl)-3*H*-quinazolin-4-one;
 6-chloro-3-(4-chloro-phenyl)-2-phenyl-3*H*-quinazolin-4-one;
 20 3-(5-chloro-pyridin-2-yl)-2-[4-(1*H*-pyrrol-2-yl)-phenyl]-3*H*-quinazolin-4-one;
 3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1*H*-pyrrol-2-yl)-phenyl]-3*H*-quinazolin-4-one;
 3-(5-chloro-pyridin-2-yl)-2-[4-(1-ethyl-1*H*-pyrrol-2-yl)-phenyl]-3*H*-quinazolin-4-one;
 25 2-{4-[1-(2-amino-ethyl)-1*H*-pyrrol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3*H*-quinazolin-4-one;
 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methylamino-ethyl)-1*H*-pyrrol-2-yl]-phenyl}-3*H*-quinazolin-4-one;
 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-ethylamino-ethyl)-1*H*-pyrrol-2-yl]-phenyl}-3*H*-quinazolin-4-one;
 30 2-{4-[1-(2-benzylamino-ethyl)-1*H*-pyrrol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3*H*-quinazolin-4-one;

- 3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-2-ylmethyl)-amino]-ethyl}-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-3-ylmethyl)-amino]-ethyl}-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 5 3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-4-ylmethyl)-amino]-ethyl}-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 2-[4-(1-benzyl-1H-pyrrol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-2-ylmethyl-1H-pyrrol-2-yl)-phenyl]-
- 10 3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-3-ylmethyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-4-ylmethyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 15 3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclohexyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(tetrahydro-pyran-4-yl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1-piperidin-4-yl-1H-pyrrol-2-yl)-phenyl]-3H-
- 20 quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(1-methyl-piperidin-4-yl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;
- 2-[4-[1-(1-acetyl-piperidin-4-yl)-1H-pyrrol-2-yl]-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- 25 3-(5-chloro-pyridin-2-yl)-2-[4-(1-isopropyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclopropyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclobutyl-1H-pyrrol-2-yl)-phenyl]-3H-
- 30 quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclopentyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

- 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(tetrahydro-furan-3-yl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(2',3',4',5'-tetrahydro-1'H-[1,3']bipyrrolyl-2-yl)-phenyl]-3H-quinazolin-4-one;
- 5 3-(5-chloro-pyridin-2-yl)-2-[4-(1'-methyl-2',3',4',5'-tetrahydro-1'H-[1,3']bipyrrolyl-2-yl)-phenyl]-3H-quinazolin-4-one;
- 2-[4-(1'-acetyl-2',3',4',5'-tetrahydro-1'H-[1,3']bipyrrolyl-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1'-methanesulfonyl-2',3',4',5'-tetrahydro-1'H-[1,3']bipyrrolyl-2-yl)-phenyl]-3H-quinazolin-4-one;
- 10 N-[2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-pyrrol-1-yl)-ethyl]-acetamide;
- 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-hydroxy-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;
- 15 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methoxy-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methoxy-1-methyl-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;
- 2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-pyrrol-1-yl)-acetamide;
- 20 2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-pyrrol-1-yl)-N-methyl-acetamide;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 25 3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1-ethyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 2-{4-[1-(2-amino-ethyl)-1H-imidazol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- 30 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methylamino-ethyl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

2-{4-[1-(2-ethylamino-ethyl)-1H-imidazol-2-yl]-phenyl}-3-pyridin-2-yl-3H-quinazolin-4-one;

2-{4-[1-(2-benzylamino-ethyl)-1H-imidazol-2-yl]-phenyl}-3-pyridin-2-yl-3H-quinazolin-4-one;

5 3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-2-ylmethyl)-amino]-ethyl}-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-3-ylmethyl)-amino]-ethyl}-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

10 2-{4-[1-(2-benzylamino-ethyl)-1H-imidazol-2-yl]-phenyl}-3-pyridin-2-yl-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-2-ylmethyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-3-ylmethyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

15 3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-4-ylmethyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclohexyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

20 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(tetrahydro-pyran-4-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-piperidin-4-yl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(1-methyl-piperidin-4-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

25 2-{4-[1-(1-acetyl-piperidin-4-yl)-1H-imidazol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-isopropyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

30 3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclopropyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclobutyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclopentyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(tetrahydro-furan-3-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

5 3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyrrolidin-3-yl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(1-methyl-pyrrolidin-3-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

10 2-{4-[1-(1-acetyl-pyrrolidin-3-yl)-1H-imidazol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(1-methanesulfonyl-pyrrolidin-3-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

N-[2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-imidazol-1-yl)-ethyl]-acetamide;

15 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-hydroxy-ethyl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methoxy-ethyl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

20 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methoxy-1-methyl-ethyl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-imidazol-1-yl)-acetamide;

2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-imidazol-1-yl)-N-methyl-acetamide;

25 2-[4-(5-amino-furan-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

2-[4-(5-aminomethyl-furan-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

30 5-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-furan-2-carboxylic acid amide;

2-{4-[5-(1-amino-1-methyl-ethyl)-furan-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

- 2-[4-(3-amino-furan-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(3-dimethylaminomethyl-furan-2-yl)-phenyl]-3H-quinazolin-4-one;
- 5 3-(5-chloro-pyridin-2-yl)-2-{4-[3-(1-dimethylamino-1-methyl-ethyl)-furan-2-yl]-phenyl}-3H-quinazolin-4-one;
- 2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-furan-3-carboxylic acid amide;
- 3-(5-chloro-pyridin-2-yl)-2-(4-oxazol-2-yl-phenyl)-3H-quinazolin-4-one;
- 10 2-[4-(5-aminomethyl-oxazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- 2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-oxazole-5-carboxylic acid amide;
- 2-[4-(4-aminomethyl-oxazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-
- 15 quinazolin-4-one;
- 2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-oxazole-4-carboxylic acid amide;
- 3-(5-chloro-pyridin-2-yl)-2-(4-thiazol-2-yl-phenyl)-3H-quinazolin-4-one;
- 2-[4-(5-aminomethyl-thiazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-
- 20 quinazolin-4-one;
- 2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-thiazole-5-carboxylic acid amide;
- 2-[4-(4-amino-thiazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- 25 N-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-thiazol-4-yl)-acetamide;
- 2-[4-(5-amino-thiazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- N-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-thiazol-5-yl)-acetamide;
- 30 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-[1,3]diazepan-1-yl)-phenyl]-3H-quinazolin-4-one;

5 2-[4-(3-amino-2-oxo-piperidin-1-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(3-dimethylamino-2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;

10 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-3-pyrrolidin-1-yl-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;

N-(1-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-2-oxo-piperidin-3-yl)-acetamide;

2-[4-(3-amino-2-oxo-pyrrolidin-1-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

15 3-(5-chloro-pyridin-2-yl)-2-[4-(2-dimethylaminomethyl-imidazol-1-yl)-phenyl]-3H-quinazolin-4-one;

1-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-1H-imidazole-2-carboxylic acid dimethylamide;

3-(5-chloro-pyridin-2-yl)-2-(4-isoxazol-5-yl-phenyl)-3H-quinazolin-4-one;

20 3-(5-chloro-pyridin-2-yl)-2-(4-oxazol-5-yl-phenyl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-(4-thiazol-5-yl-phenyl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(3H-[1,2,3]triazol-4-yl)-phenyl]-3H-quinazolin-4-one;

25 3-(5-chloro-pyridin-2-yl)-2-[4-(5-methyl-4H-[1,2,4]triazol-3-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(5-methyl-[1,3,4]thiadiazol-2-yl)-phenyl]-3H-quinazolin-4-one;

2-biphenyl-4-yl-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

2-(2'-amino-biphenyl-4-yl)-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

30 3-(5-chloro-pyridin-2-yl)-2-(2'-dimethylamino-biphenyl-4-yl)-3H-quinazolin-4-one;

2-(2'-aminomethyl-biphenyl-4-yl)-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

- 3-(5-chloro-pyridin-2-yl)-2-(2'-dimethylaminomethyl-biphenyl-4-yl)-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-(4-pyridin-2-yl-phenyl)-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-(4-pyridin-3-yl-phenyl)-3H-quinazolin-4-one;
- 5 3-(5-chloro-pyridin-2-yl)-2-(4-pyridin-4-yl-phenyl)-3H-quinazolin-4-one;
- 2-[4-(2-amino-pyridin-3-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- 2-[4-(2-aminomethyl-pyridin-3-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;
- 10 3-(5-chloro-pyridin-2-yl)-2-[4-(2-dimethylaminomethyl-pyridin-3-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 15 6-chloro-3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-6-fluoro-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 20 6-bromo-3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-4-oxo-3,4-dihydro-quinazoline-6-carbonitrile;
- 3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-6-methoxy-3H-quinazolin-4-one;
- 25 3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-4-oxo-3,4-dihydro-quinazoline-6-carboxylic acid amide;
- 6-chloro-3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 30 3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-6-methoxy-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-6-fluoro-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;

- 3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-4-oxo-3,4-dihydro-quinazoline-6-carbonitrile;
- 3-(4-chloro-phenyl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 5 2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3-(4-methoxy-phenyl)-3H-quinazolin-4-one;
- 3-(3-chloro-phenyl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 10 2-fluoro-5-{2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-4-oxo-4H-quinazolin-3-yl}-benzonitrile;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-pyrrolidin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(4-chloro-phenyl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(4-methoxy-phenyl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 15 3-(3-chloro-phenyl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(3-chloro-phenyl)-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(4-chloro-phenyl)-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 20 3-(5-chloro-pyridin-2-yl)-6-methoxy-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-dimethylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-6-methoxy-3H-quinazolin-4-one;
- 25 6-chloro-3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-dimethylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;
- 6-chloro-3-(4-chloro-phenyl)-2-{4-[1-(2-dimethylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;
- 3-(5-chloro-pyridin-2-yl)-6-methoxy-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 30 6-chloro-3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;
- 6-chloro-3-(4-chloro-phenyl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;

6-chloro-3-(4-methoxy-phenyl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;

6-chloro-3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-quinazolin-4-one;

5 3-(5-chloro-pyridin-2-yl)-6-methoxy-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-6-methoxy-3H-quinazolin-4-one;

10 3-(4-chloro-phenyl)-2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-6-methoxy-3H-quinazolin-4-one;

6-chloro-3-(4-chloro-phenyl)-2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-3H-quinazolin-4-one;

6-bromo-3-(4-chloro-phenyl)-2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-3H-quinazolin-4-one;

15 2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-3-(4-methoxy-phenyl)-6-methyl-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;

20 3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-[1-(2-methylamino-ethyl)-1H-pyrrol-2-yl]-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-pyrido[4,3-d]pyrimidin-4-one;

25 3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-pyrido[4,3-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-pyrido[3,4-d]pyrimidin-4-one;

30 3-(5-chloro-pyridin-2-yl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-thieno[3,4-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-imidazol-2-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;

- 3-(5-chloro-pyridin-2-yl)-2-[4-(1H-imidazol-2-yl)-phenyl]-3H-thieno[3,4-d]pyrimidin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1H-imidazol-2-yl)-phenyl]-3H-thieno[3,4-d]pyrimidin-4-one;
- 5 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-pyrido[4,3-d]pyrimidin-4-one;
- 10 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-pyrido[3,4-d]pyrimidin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-thieno[3,4-d]pyrimidin-4-one;
- 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;
- 15 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-pyrido[4,3-d]pyrimidin-4-one;
- 6-(5-chloro-pyridin-2-yl)-5-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-6H-[1,2,5]oxadiazolo[3,4-d]pyrimidin-7-one;
- 6-(5-chloro-pyridin-2-yl)-5-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-6H-isoxazolo[4,3-d]pyrimidin-7-one;
- 20 6-(5-chloro-pyridin-2-yl)-5-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-6H-thiazolo[5,4-d]pyrimidin-7-one;
- 1-(5-chloro-pyridin-2-yl)-7-methyl-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-1,7-dihydro-purin-6-one;
- 25 3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-thieno[3,2-d]pyrimidin-4-one;
- 3-(5-chloro-pyridin-2-yl)-5-methyl-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3,5-dihydro-pyrrolo[3,2-d]pyrimidin-4-one;
- 6-(5-chloro-pyridin-2-yl)-1-methyl-5-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-1,6-dihydro-pyrazolo[4,3-d]pyrimidin-7-one;
- 30 3-(5-chloro-pyridin-2-yl)-2-[4-(3-dimethylamino-2-oxo-piperidin-1-yl)-phenyl]-3H-furo[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(3-dimethylamino-2-oxo-piperidin-1-yl)-phenyl]-5-methyl-3,5-dihydro-pyrrolo[3,2-d]pyrimidin-4-one; and,

6-(5-chloro-pyridin-2-yl)-5-[4-(3-dimethylamino-2-oxo-piperidin-1-yl)-phenyl]-1-methyl-1,6-dihydro-pyrazolo[4,3-d]pyrimidin-7-one;

5 or a pharmaceutically acceptable salt form thereof.

9. A method for treating a thromboembolic disorder, comprising:
administering to a patient in need thereof a therapeutically effective amount of a
compound of Claim 1 or a pharmaceutically acceptable salt thereof.

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10. A method according to Claim 9, wherein the thromboembolic disorder is
selected from the group consisting of arterial cardiovascular thromboembolic
disorders, venous cardiovascular thromboembolic disorders, and thromboembolic
disorders in the chambers of the heart.

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11. A method according to Claim 9, wherein the thromboembolic disorder is
selected from unstable angina, an acute coronary syndrome, first myocardial
infarction, recurrent myocardial infarction, ischemic sudden death, transient ischemic
attack, stroke, atherosclerosis, peripheral occlusive arterial disease, venous
20 thrombosis, deep vein thrombosis, thrombophlebitis, arterial embolism, coronary
arterial thrombosis, cerebral arterial thrombosis, cerebral embolism, kidney embolism,
pulmonary embolism, and thrombosis resulting from (a) prosthetic valves or other
implants, (b) indwelling catheters, (c) stents, (d) cardiopulmonary bypass, (e)
hemodialysis, and (f) other procedures in which blood is exposed to an artificial
25 surface that promotes thrombosis.

12. A method of treating a patient in need of thromboembolic disorder
treatment, comprising: administering a compound of Claim 1 or a pharmaceutically
acceptable salt thereof in an amount effective to treat a thromboembolic disorder.

30

13. A method, comprising: administering a compound of Claim 1 or a
pharmaceutically acceptable salt thereof in an amount effective to treat a
thromboembolic disorder.